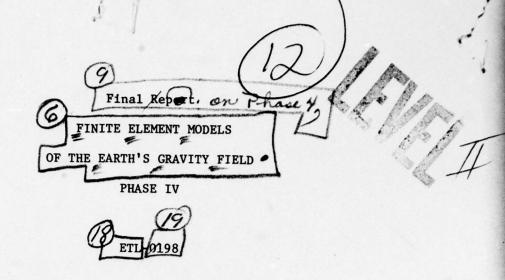


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U.S. Army Engineer Topographic Laboratories Fort Belvoir, Virginia 22060

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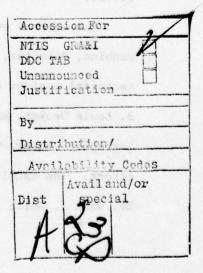
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### 1.0 Summary

This research and development effort has produced the following results:

- \* Software suitable for the construction of finite element gravity disturbance fields.
- \* Software suitable for calculation of the gravity disturbances within finite element fields:
  - using Chebyshev polynomials
  - using Orthonormal polynomials
- \* Example finite element gravity disturbance fields have been developed and tested. Relevant comparisons between the Chebyshev polynomials approach and the Orthonormal polynomials approach have been made.
- \* Optimizations have been incorporated to increase computational efficiency and to determine a model for the geopotential, using Orthogonal polynomials.



#### 2.0 Preface

This report constitutes the final report under Contract No. DAA 70-78-C-0072 performed by the Virginia Polytechnic Institute and State
University for the U. S. Army Engineer Topographic Laboratories, Fort
Belvoir, Virginia, under the sponsorship of the Defense Mapping Agency.

The current effort is primarily concerned with the investigation of new choices of computationally more efficient basis functions which can be used to model the local fine structure of the gravity vector. This development, to date, has produced two separate versions of software systems. This report documents each of these systems as well as their theoretical background. Comparisons are made based on the result of several test cases. Also included are detailed discussions of the software and standard print-outs of each version. In addition, the programs contain enough detailed comment statements to allow the know-ledgeable user to use and modify particular sections to suit his or her own needs. Finally, the report also indicates a possible approach to determine a model for the geopotential itself, which at the same time would increase the computational efficiency of determining the gravity vector.

The authors very much appreciate the technical liaison of Mr. L. A. Gambino, Director, Computer Sciences Laboratory (USAETL), who served as technical monitor of this work. Discussions with Dr. R. W. Ballew, Dr. B. Louis Decker, and Mr. H. W. Howard (all of the Defense Mapping Agency, Aerospace Center) were also helpful in defining the objectives of our effort. The excellent programming support of Mr. John T. Saunders and Mr. John J. Smith is gratefully acknowledged.

#### 3.0 Introduction

The central theme in the development of the present software systems is the finite element approach to piecewise approximation. The finite element concept may be considered as an extension to three dimensions of one dimensional piecewise approximation techniques. Qualitatively, some quantity that is functionally related in a "complicated" fashion to position over a "large" interval may be replaced for much "smaller" intervals of position by much "less complicated" approximating functions. Thus the complicated function, which will normally be "more expensive" to evaluate, will have been replaced by a set of locally valid functions which will normally be "less expensive" to evaluate.

The idea may easily be extended to three dimensions. A quantity which is functionally related to position over some large region may be replaced by a set of approximations, each of which is valid only in a small local volume; the set of local volumes spanning the large region of interest. The local approximations are typically much easier (faster computationally) to evaluate. The finite element concept may be applied to any of a variety of modeling problems; this investigation centered specifically on fine structure gravity modeling.

Previously, earth-fixed spherical coordinate based finite element fields have been successfully used to replace globally valid spherical harmonic representations of the geopotential and its derivatives. But because the model (in the present study) used to simulate gravity disturbance data (Model 310) consists of a set of point masses placed in relation to the Geodetic Reference Surface of 1967, ellipsoidal coordinates (H,  $\lambda$ ,  $\phi$  - see Figure 1) were used in the finite element modeling process.

Finite element fields may be defined by ellipsoidal coordinates as shown in Figure 2. The region to be modeled is defined by a set of maximum bounds ( $H_{max}$ ,  $\lambda_{max}$ ,  $\phi_{max}$ ) and a set of minimum bounds ( $H_{min}$ ,  $\lambda_{min}$ ,  $\phi_{min}$ ). The region is then divided into a set of smaller volumes, or finite element cells, where each cell has dimensions  $H_{cell}$ ,  $\lambda_{cell}$ ,  $\phi_{cell}$ . Thus the 300 km  $\times$  10° region above the reference ellipsoid, altitude 0 - 300 km,

altitude 0 - 300 km, longitude  $70^{\circ}\text{E} - 80^{\circ}\text{E}$ , and latitude  $25^{\circ}\text{S} - 35^{\circ}\text{S}$ ,

could be modeled by one hundred (100) cells each 300 km  $\times$  1°  $\times$  1° or four (4) cells each 300 km  $\times$  5°  $\times$  5° or one cell 300 km  $\times$  10°  $\times$  10°, etc.

The approximation valid in any one cell is generated independently of the approximation for any other cell. A set of gravity disturbance "observations" are produced for each cell directly from Model 310 and then approximated for only that cell in a least squares fitting process. The approximating models thus generated are then stored on a high speed rotating mass storage device for later use in the computation of the approximate gravity disturbance in any finite element cell of the region modeled.

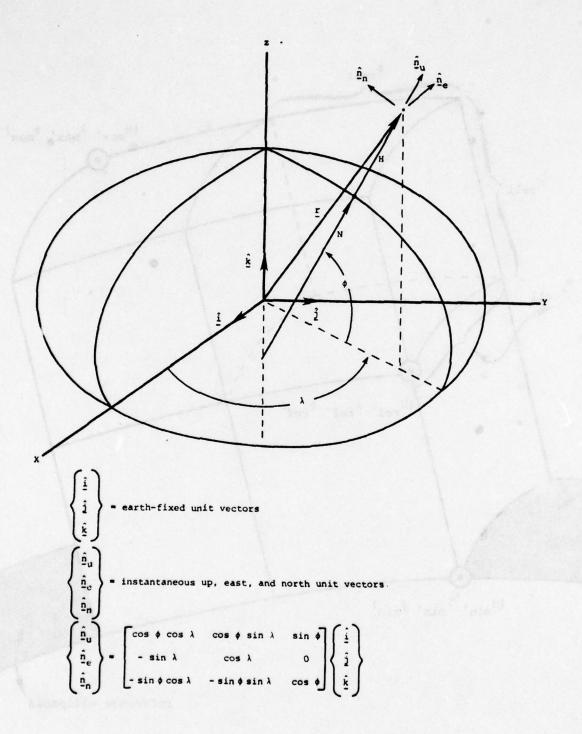


Figure 1 Earth-Fixed Rectangular and Ellipsoidal Coordinate Systems.

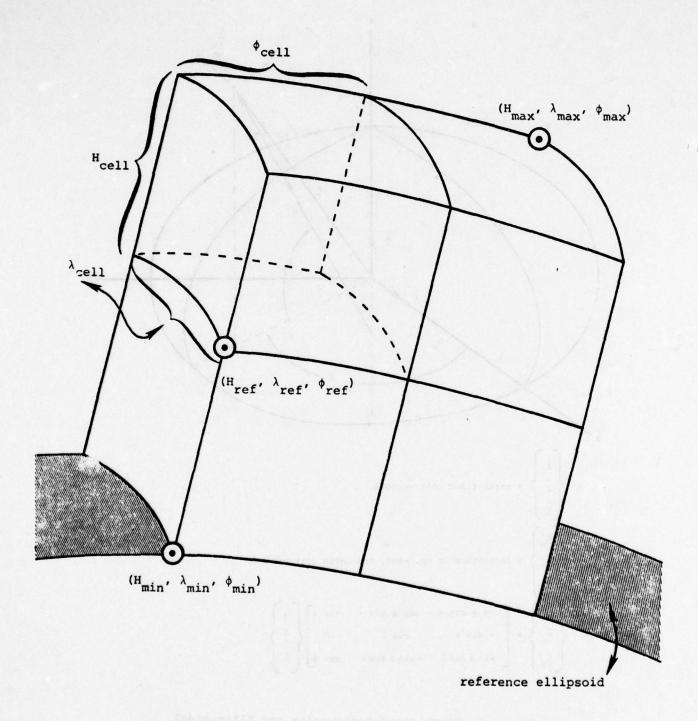


Figure 2 Ellipsoidal Finite Element Field.

## 4.0 Finite Element Formulation

Define the gravity disturbance, og:

$$\underline{g}_{\text{actual}} = \underline{g}_{\text{reference}} + \delta \underline{g}$$
 (1)

$$\delta \underline{g} = \begin{cases} \delta g_{up} \\ \delta g_{east} \end{cases} = \underline{g}_{actual} - \underline{g}_{reference}$$
 (2)

Now, replace the components of  $\delta \underline{g}$  with locally valid polynomial approximations,  $\delta \underline{G}$ , of the form,

$$\delta \underline{G} = \begin{cases}
\delta G_{up} \\
\delta G_{east} \\
\delta G_{north}
\end{cases} = \begin{cases}
\sum_{n=0}^{N} \sum_{i=0}^{n} \sum_{j=0}^{C_{up}} \begin{cases} C_{up_{ijk}} \\
C_{east_{ijk}} \\
C_{north_{ijk}}
\end{cases} f_{ijk}(X_1, X_2, X_3)$$
(3)

where,

N = NORDER = the highest order polynomial,

$$k = n - i - j$$

$$X_1 = (H - H_{ref})/H_{cell}$$

$$X_2 = (\lambda - \lambda_{ref})/\lambda_{cell}$$
 (non-dimensional coordinates) (4)

$$X_3 = (\phi - \phi_{ref})/\phi_{cell}$$

 $f_{ijk}$  = chosen polynomial basic functions, and

 $c_{up}^{\phantom{\dagger}}_{ijk}$ ,  $c_{east}^{\phantom{\dagger}}_{ijk}$ , and  $c_{north}^{\phantom{\dagger}}_{ijk}$  are constant coefficients determined via

least square fits so that  $\sum_{i=1}^{m} (\delta G_{up} - \delta g_{up})_{i}^{2}$ ,  $\sum_{i=1}^{m} (\delta G_{east} - \delta g_{east})_{i}^{2}$ ,

and  $\sum_{i=1}^{m} (\delta G_{north} - \delta g_{north})_{i}^{2}$  are minimized for some local volume of

space (i.e. the cell being modeled).

Thus, to determine the three sets of constant coefficients,  $C_{\rm up}$ ,  $C_{\rm east}$ , and  $C_{\rm north}$ , for each cell, the following linear least squares problems must be solved,

$$||\delta\underline{G}_{up} - \delta\underline{g}_{up}|| = ||\underline{AC}_{up} - \delta\underline{g}_{up}|| = \min.$$

$$||\delta\underline{G}_{east} - \delta\underline{g}_{east}|| = ||\underline{AC}_{east} - \delta\underline{g}_{east}|| = \min.$$

$$||\delta\underline{G}_{north} - \delta\underline{g}_{north}|| = ||\underline{AC}_{north} - \delta\underline{g}_{north}|| = \min.$$
(5)

where the coefficient matrix, A, has the form,

$$A = \begin{bmatrix} f_{000}^{(1)} & f_{001}^{(1)} & f_{010}^{(1)} & \dots & f_{N00}^{(1)} \\ f_{000}^{(2)} & f_{001}^{(2)} & f_{010}^{(2)} & \dots & f_{N00}^{(1)} \\ & & & & & & \\ f_{000}^{(m)} & f_{001}^{(m)} & f_{010}^{(m)} & \dots & f_{N00}^{(m)} \end{bmatrix}$$
(6)

in which m is the number of observations in the cell and  $f_{ijk}(\ell)$  represents  $f_{ijk}(X_1, X_2, X_3)$  corresponding to the  $\ell$ th observation.

Furthermore, the three coefficient vectors,  $\underline{C}_{up}$ ,  $\underline{C}_{east}$ , and  $\underline{C}_{north}$ , have the form

$$\frac{C_{up}_{000}}{C_{up}_{001}} = \begin{pmatrix} C_{east}_{000} \\ C_{up}_{001} \\ C_{up}_{010} \\ C_{up}_{100} \\ C_{up}_{100} \\ C_{up}_{002} \\ C_{up}_{001} \\ C_{up}_{002} \\ C_{up}_{001} \\ \vdots \\ C_{up}_{no0} \end{pmatrix}, \quad \frac{C_{east}}{C_{east}_{000}} = \begin{pmatrix} C_{north}_{000} \\ C_{north}_{010} \\ C_{north}_{100} \\ C_{north}_{100} \\ C_{north}_{002} \\ C_{north}_{011} \\ \vdots \\ C_{north}_{north}_{001} \\ \vdots \\ C_{north}_{north}_{000} \end{pmatrix} \tag{7}$$

where

N = NORDER

and the three sets of gravity disturbance observations,  $\delta g_{up}$ ,  $\delta g_{east}$ , and  $\delta g_{north}$ , are

$$\delta \mathbf{g}_{up} = \begin{cases} \delta \mathbf{g}_{up}(\mathbf{X}_{1_{1}}, \mathbf{X}_{2_{1}}, \mathbf{X}_{3_{1}}) \\ \delta \mathbf{g}_{up}(\mathbf{X}_{1_{2}}, \mathbf{X}_{2_{2}}, \mathbf{X}_{3_{2}}) \\ \delta \mathbf{g}_{up}(\mathbf{X}_{1_{3}}, \mathbf{X}_{2_{3}}, \mathbf{X}_{3_{3}}) \\ \vdots \\ \delta \mathbf{g}_{up}(\mathbf{X}_{1_{m}}, \mathbf{X}_{2_{m}}, \mathbf{X}_{3_{m}}) \end{cases}$$

$$\delta \mathbf{g}_{east} = \begin{cases} \delta \mathbf{g}_{east}(\mathbf{X}_{1_{1}}, \mathbf{X}_{2_{1}}, \mathbf{X}_{3_{1}}) \\ \delta \mathbf{g}_{east}(\mathbf{X}_{1_{2}}, \mathbf{X}_{2_{2}}, \mathbf{X}_{3_{2}}) \\ \delta \mathbf{g}_{east}(\mathbf{X}_{1_{3}}, \mathbf{X}_{2_{3}}, \mathbf{X}_{1_{3}}) \\ \vdots \\ \delta \mathbf{g}_{east}(\mathbf{X}_{1_{m}}, \mathbf{X}_{2_{m}}, \mathbf{X}_{3_{m}}) \end{cases}$$

$$\delta g_{north} = \begin{cases} \delta g_{north}(X_{1_{1}}, X_{2_{1}}, X_{3_{1}}) \\ \delta g_{north}(X_{1_{2}}, X_{2_{2}}, X_{3_{2}}) \\ \delta g_{north}(X_{1_{3}}, X_{2_{3}}, X_{3_{3}}) \\ \vdots \\ \delta g_{north}(X_{1_{m}}, X_{2_{m}}, X_{3_{m}}) \end{cases}$$
(8)

where  $\{(X_1, X_2, X_3), i = 1, 2, ..., m\}$  = a specified set of  $(H_i, \lambda_i, \phi_i)$  coordinates - usually a uniform observation grid in any one cell. The gravity disturbance observations are determined directly from Model 310 evaluations on the observation grid in each cell.

Following the standard least squares approximation procedure we obtain, for the "up-direction"

$$\underline{\mathbf{C}}_{\mathbf{u}\mathbf{p}} = \mathbf{B} \, \delta \underline{\mathbf{g}}_{\mathbf{u}\mathbf{p}} \tag{9}$$

where

$$B = (A^{T} A)^{-1}A^{T}$$
 (10)

Similar results are valid in the "east- and north-directions".

## 5.0 Chebyshev Polynomials Finite Element Models

The choice of basis functions in Eq. (3) is still open. Several sets of basis functions have been used to date. Polynomial sets are of particular interest because of the ease with which they can be manipulated. The most obvious set is given by

Another set consists of products of shifted Chebyshev polynomials

Chebyshev polynomials,  $t_n(X_1)$ , of order n may be generated as follows:

For  $n \ge 2$ , a recursion formula may be used:

$$t_n(X) = 2X t_{n-1}(X) - t_{n-2}(X), -1 \le X \le 1$$
 (13)

Shifted Chebyshev polynomials,  $T_n(\overline{X})$  may be computed by substituting

$$X = 2\overline{X} - 1$$

so that,

$$T_n(\bar{X}) = t_n(X) = t_n(2\bar{X} - 1), \qquad 0 \le \bar{X} \le 1$$
 (14)

To generate shifted Chebyshev polynomials a set of X's with values between 0 and 1 are needed. The difference between some point  $(H,\lambda,\phi)$  in a cell and a reference point  $(H_{ref},\lambda_{ref},\phi_{ref})$  (the "lowermost corner"

of the cell) is divided by the cell's dimensions,  $H_{cell}$ ,  $\lambda_{cell}$ ,  $\phi_{cell}$ , as in (4) to obtain a set of non-dimensional coordinates,  $X_1$ ,  $X_2$ ,  $X_3$ , for each cell. The non-dimensional coordinates may then be used to generate the shifted Chebyshev polynomials.

#### 5.1 Discussion

Certain aspects of the process by which finite element fields are created may be capitalized on under special circumstances. When it is known beforehand that uniformly gridded gravity disturbance observations will be available, only one least squares matrix (A matrix) need be generated. This is true because the set of non-dimensional coordinates,  $\{(X_{1_i}, X_{2_i}, X_{3_i}), i=1, 2, \ldots, m\} \text{ of one cell's observation grid will be the same as every other cell's non-dimensional observation grid coordinates when the observation grid pattern is set for all cells. If the same number of observations are made in each cell and the positions of the observations points are the same relative to the reference point (<math>H_{\text{ref}}$ ,  $h_{\text{ref}}$ ,  $h_{\text{ref}}$ ) in each cell, then all cells will have the same A matrix. The A matrix, after all, consists only of products of the locally evaluated polynomials, which are functions only of the non-dimensional coordinates of the observation points.

This all means that only one A-matrix need be generated for the first cell of a finite element field; but the gravity observations must be generated for each cell. A reduction of the A-matrix to upper triangular form may be performed for the least squares fitting process, after which each set of observations may be similarly reduced and then back-substituted to produce a set of coefficients for each cell. The A-

matrix, again, needs to be reduced only once.

It should be noted that this method of finite element field generation does not rely on Model 310 or any other particular gravity disturbance model. Model 310 could be replaced by any other process that is capable of producing the gravity disturbance observations at the observation grid points of each cell.

Once a finite element field has been created, the Chebyshev polynomial gravity disturbance approximation  $\delta \underline{G}$  may be computed by evaluating (3). The subscripts ijk always follow the same pattern for a given NORDER. This means that a "current coefficient number",  $\ell$ , may be attached to each group of subscripts ijk and that the total number of coefficients, NC, necessary to model any one component of the gravity disturbance in one cell will be constant for a given NORDER. Table 1 shows these relationships for various NORDER's.

Since we can precompute and save the subscripts, ijk, in the appropriate pattern such that they are functions of only the current coefficient number,  $\ell$ , (3) can be written,

$$\delta \underline{G} = \begin{cases} \delta G_{up} \\ \delta G_{east} \\ \delta G_{north} \end{cases} = \begin{cases} \sum_{\ell=1}^{NC} \begin{cases} C_{up_{\ell}} \\ C_{east_{\ell}} \\ C_{north_{\ell}} \end{cases} T_{i}(X_{1}) T_{j}(X_{2}) T_{k}(X_{3})$$
(15)

where,

NC = the total number of coefficients for one component of the gravity disturbance, as above

ijk = precomputed and saved as functions of  $\ell$ 

 $X_1, X_2, X_3$  = functions of H, $\lambda$ , $\phi$  and the constants  $H_{ref}$ ,  $\lambda_{ref}$ ,  $\phi_{ref}$ ,  $H_{cell}$ ,  $\lambda_{cell}$ ,  $\phi_{cell}$  as in (4) above. The total number of summations in (15) is the same as in (3) but the amount of bookkeeping has been cut down tremendously by doing business in this fashion. Computational savings of up to 50% were obtained by using (15) in lieu of (3) for evaluation purposes. Table 1 was also used in the generation of the Amatrix, but the saving here is not as significant due to the previously discussed fact that the A-matrix is only generated once.

NORDER	ijk kastra	<u>L</u> vd	here z <u>NC</u>
atolika a un Oone ya	0 0 0	ysolalals	e glasifus sa webs
of oldison 21 i) as	0 0 1	all gunyio	g lawronodits fo the
	0 1 0	3 10 11	
to singeons rebut	100		eldraeoq of lllw sl
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	0 1 1	6	te testaseés for us
	0 2 0	7	or an 70, 10% or an
	101	8	
	1 1 0	9	
	2 0 0	10	10
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	0 1 2	12	
	:::	•	
	3 0 0	20	20
•	• • • /		•
	1121.610 8 8.811	130.0	1.0)
N	0 0 N		
	0 1 N-1	2 9,102 (01)	
	:::	1.11	
	N O O	NC	$\begin{array}{ccc} N+1 & n \\ \sum & \sum m \\ n=1 & m=1 \end{array}$

-

Table 1 Subscript Patterns for Various NORDER's.

#### 6.0. Orthonormal Polynomials Finite Element Models

As demonstrated by the numerical results, Chebyshev polynomials yield an entirely satisfactory solution. However, by choosing a judicious set of orthonormal polynomials as basis functions, it is possible to avoid the computation of the matrix inverse in Eq. (10). In addition, it will be possible to compute the coefficients  $c_{ijk}$  independently of each other. This is especially convenient when the order of approximation is increased for no need exists to recompute the least squares operator B in Eq. (10), or any lower order coefficients.

Let us introduce the following basis functions

Where  $f_i$ ,  $f_j$ ,  $f_k$  represent (to be specified) one-dimensional polynomials in  $X_1$ ,  $X_2$ , and  $X_3$  respectively.

Using Eq. (6), it is not difficult to show that the matrix  $\mathbf{A}^{T}\mathbf{A}$  has elements of the form

$$\sum_{\ell=1}^{m} f_{ijk}(\ell) f_{\alpha\beta\gamma}(\ell) ; i,j,k = 0,1,...,N$$

$$\alpha,\beta,\gamma = 0,1,...,N$$
(17)

and introducing Eq. (16) into Eq. (17) yields

$$\sum_{\ell_{1}=1}^{m_{1}} f_{i}(\ell_{1}) f_{\alpha}(\ell_{1}) \sum_{\ell_{2}=1}^{m_{2}} f_{j}(\ell_{2}) f_{\beta}(\ell_{2}) \sum_{\ell_{3}=1}^{m_{3}} f_{k}(\ell_{3}) f_{\gamma}(\ell_{3})$$
(18)

where  $m_i$  (i = 1,2,3) represents the number of observations in the  $X_i$ direction, and m is the total number of observations in the cell.

Next, let us assume that the set  $f_i(X_1)$  can be constructed orthonormal over the discrete domain of interest. Hence,

$$\sum_{\ell_1=1}^{m_1} f_i(\ell_1) f_{\alpha}(\ell_1) = \delta_{i\alpha}$$
(19)

Similar expressions can be written for the sets  $f_j(X_2)$  and  $f_k(X_3)$ . Substituting Eq. (19) into Eq. (18) yields

$$^{\delta}_{\mathbf{i}\alpha}{^{\delta}_{\mathbf{j}}\beta}{^{\delta}_{\mathbf{k}}\gamma} \tag{20}$$

and because (20) is unity if and only if  $i=\alpha$ ,  $j=\beta$ ,  $k=\gamma$ , we conclude that the matrix  $A^TA$  reduces to the unit matrix.

The problem is now reduced to the construction of an orthonormal set of one-dimensional polynomials  $f_i(X)$  (i = 0,1,...,N).

Any orthogonal set of polynomials has the following three-term recurrence relation[3]

$$\frac{h_{i}}{h_{i+1}} f_{i+1}(x) = x f_{i}(x) - \frac{g_{i}h_{i-1}}{g_{i-1}h_{i}} f_{i-1}(x) - B_{i}f_{i}(x)$$
 (21)

where  $h_i$  stands for the leading coefficient in  $f_i(x)$  and  $B_i$  is an unknown constant, and where we use the inner product notation

$$\mathbf{g_i} = \langle \mathbf{f_i}, \mathbf{f_i} \rangle = \int_{\mathbf{D}} \mathbf{f_i^2}(\mathbf{x}) d\mathbf{x}$$
 (22)

and where D is the domain of interest. If we choose the following non-dimensional coordinates

$$x_1 = \frac{2(H-H_m)}{H_{cell}}$$

$$X_2 = \frac{2(\lambda - \lambda_m)}{\lambda_{cell}}$$
 (23)

$$x_3 = \frac{2(\phi - \phi_m)}{\phi_{cell}}$$

then, the domain D is (-1,1) which enables us to conclude that the functions  $f_0 = 1$  and  $f_1 = x$  are orthogonal over D. Moreover, using the Gram-Schmidt procedure it can be shown that  $f_{2k}(k=0,1,\ldots)$  is even while  $f_{2k+1}(k=0,1,\ldots)$  is odd. From Eq. (21) it then follows that all  $B_i$  must be equal to zero. Finally, because  $h_0 = h_1 = 1$ , we also must have  $h_i = 1$  ( $i=2,3,\ldots$ ). Relation (21) then reduces to

$$f_{i+1}(x) = xf_i - \frac{\langle f_i, f_i \rangle}{\langle f_{i-1}, f_{i-1} \rangle} f_{i-1}(x)$$
 (24)

which, together with  $f_0 = 1$  and  $f_1 = x$ , yields a recurrence relation for a set of orthogonal polynomials. Normalization is accomplished by

$$\hat{f}_{i} = \frac{f_{i}}{\langle f_{i}, f_{i} \rangle^{1/2}} \quad i = 0, 1, ..., N$$
 (25)

where  $\hat{f}_{i}$  represents the normalized polynomial. For a discrete observation grid we replace Eq. (22) by

$$\langle f_i, f_i \rangle = \sum_{\ell_1=1}^{m_1} f_i^2(\ell_1)$$
 (26)

It is also necessary that the local coordinates of the grid points in a particular direction add up to zero. Furthermore, we must have that the number of grid points in a particular direction is at least N+1.

Specifically Table 2 gives the orthonormal polynomials  $f_i$  (i = 0,1,2,3) for the grid points (-1,-1/3,1/3,1).

$$\hat{f}_0 = .500000$$

$$\hat{f}_1 = .670820X$$

$$\hat{f}_2 = 1.125000X^2 - .625000$$

$$\hat{f}_3 = 2.51558X^3 - 2.29197X$$

Table 2 Orthonormal Polynomials

Introducing the polynomials (25) into Eq. (16) and into Eq. (6) reduces Eq. (10) to

$$B = A^{T}$$
 (27)

so that, from Eq. (9) it follows that

$$\underline{\underline{C}}_{up} = A^{T} \delta \underline{g}_{up} \tag{28}$$

It is then possible to write an explicit expression for each coefficient  $\mathbf{c}_{\mathbf{ijk}}$  as follows

$$c_{ijk_{up}} = \sum_{\ell=1}^{m} f_{ijk}(\ell) \delta g_{up}(\ell)$$
 (29)

Note that each coefficient  $c_{ijk}$  can be computed independently and no need exists to recompute B when the value of N changes. This feature is very valuable when, for example, a prototype gravity model of a launch-region is desired but it is difficult to specify the order N required for an accurate representation; also note that N could be variable depending upon the frequency and amplitude of local gravity anomalies. Also, in case of a global gravity model, we can use a truncation of a spherical harmonic series for  $\delta g_{up}(\ell)$  and compute

$$c_{ijk_{up}} = \int_{D} f_{ijk} \delta g_{up} dD$$
 (30)

by carring out analytical integrations. This reduces Eq. (3) to a truncated Fourier series. Because a continuous spectrum of "measurements" is used, we can expect more accurate approximations.

#### 6.1 Discussion

The recursive relationship of the orthonormal polynomials is very similar to the relationship previously used with Chebyshev polynomials. This formulation however does contain two additional division operations as well as a square root evaluation which must be performed on each polynomial. If one division and the square root values are saved during the preliminary calculation (SUBROUTINE ORTHO) and passed on for the actual evaluation, (SUBROUTINE MULT) a time savings can be realized at the small expense of a few (42) additional "words" of storage.

Note that even with this saving, the original Chebyshev polynomials can still be evaluated slightly faster, the difference only being the execution of a single division operation. However, when the advantages noted earlier are taken into account, the choice of orthonormal polynomials may be more beneficial, especially considering the effort that is involved with several different "set-ups" of variable polynomial orders and/or observation grid patterns where the Chebyshev version would require extensive computational time to process the least squares matrix and evaluate the coefficients. Indeed, the orthonormal approach does not require evaluation of  $(\mathbf{A}^{\mathrm{T}}\mathbf{A})^{-1}$  and in addition, every coefficient  $\mathbf{c}_{ijk}$  is independently computed.

#### 7.0 Discussion of the Software

A variety of Fortran programs and subroutines were written during the course of this research effort. The software developed by Junkins and Saunders<sup>2</sup> was converted from its original CDC form to run on the VPI&SU IBM System 360 Model with 158 Processors. After correctly reproducing their test results, the software was modified to incorporate some programming efficiencies. From this starting point, new programs were developed to implement the orthonormal approach as well as test various additional approaches.

In general, the final software package can be divided into three separate stand-alone modules, regardless of the basic functions chosen.

Section I: A Mass Model 310 Initialization Section.

Section II: A Least Squares Reduction and Coefficient Determination Section.

Section III: A Test Section.

Section I, A Mass Model 310 Initialization Section, consists of a single fortran program, MASPOS, and its coded input data. This job step is identical for all the versions of the software developed so far and it will be required by any new version which makes use of Mass Model 310 to supply simulated gravity disturbances for use as observations or as a control to compare the modeling equation's calculations against. Clearly, this module can be replaced by input software if other sources (e.g., measurements) of gravity disturbances are available.

This program presently inputs coded values of the 1080 point masses

of Mass Model 310. Their location's are converted from the geodetic coordinates  $(H,\lambda,\phi)$  to rectangular coordinates (x,y,z) and the mass codes are converted to their actual values. This information is then written onto a sequential access file (FILE 2), for use by SUBROUTINE PTMASS, which is described elsewhere. More specifically, PROGRAM MASPOS creates an unformatted sequential file, FILE 2, which contains the precomputed products of the gravitational constant and each point mass of Model 310 and the precomputed geocentric rectangular coordinates of each point mass. PROGRAM MASPOS requires as input coded information about Model 310 in the form of 30 cards. Each card represents 36 mass points in a grid row of equal latitude where the mass points are 50 longitude minutes apart. The mass points should be listed on each card by increasing longitude and the cards should be arranged in order of decreasing latitude. PROGRAM MASPOS reads only the coded multiplication factors for each of the 36 mass points of the grid row, where the coded factors indicate the mass of the respective mass points,  $1 = -1.\times10^{19}$ grams, 2 = 0. grams, and  $3 = +1.\times10^{19}$  grams. The coded factors are located in columns 25-60 of each data card. All other data card columns are ignored (for information only, columns 1-9 indicate south latitude in arc minutes of the grid row and columns 10-18 indicate east longitude in arc minutes of the west-most mass point in the row).

Section II, A Least Squares Reduction and Coefficient Determination

Section, is where the actual modeling equation coefficients are calculated.

Since the coefficients are determined by the choice of the basis functions and the choice of basis functions is a primary issue under study, each version of the software for Section II differs in several important aspects while the fundamental sections of the programs, such as establishment of

the finite element cells and grid pattern and determination of the gravity disturbances will be the same.

The main program for Section IIA, LOCALG, performs all the necessary preliminary calculations that are required by each version. Using card input, the finite element field is established and broken down into finite element cells and the observation grid pattern is also established. The critical field information is then written onto a storage file (FILE 3) for later use by PROGRAM FINEG. More specifically, PROGRAM LOCALG generates the A-matrix for a "typical" finite element cell and a set of gravity disturbance observations for each cell of a finite element gravity disturbance field. The gravity disturbance observations are evaluations of SUBROUTINE PTMASS on the uniform observation grid in each cell. The user defines a finite element gravity disturbance field by providing the following items as inputs to PROGRAM LOCALG:

- The integral order of the locally valid polynomial approximating functions, NORDER, (0 < NORDER < 6).</li>
- 2. The observation grid pattern (same for all cells) to include the number of observations in the up, eastern and northern directions, MOBSU, MOBSE, MOBSN. The total number of observations (MOBSU × MOBSE × MOBSN) should be at least three times the number of coefficients, NC (see Table 1), in the approximating model. Furthermore, the number of

observations in each direction should be one more than NORDER (MOBSU = MOBSE = MOBSN = NORDER + 1). The maximum number of observations allowed by the present program dimensions is 343 (a  $7 \times 7 \times 7$  observation grid). Non-uniform sample grids (e.g.  $5 \times 7 \times 7$ ,  $6 \times 5 \times 4$ , etc.) have been found to produce unpredictable results - often an unacceptable fit.

- 3. The size of each finite element cell in H,λ,φ must be specified, HCELL, ALCELL, APCELL. HCELL is cell size in H (meters), (HCELL ≤ [HMAX HMIN]). ALCELL is cell size in λ (degrees), (ALCELL ≤ [ALMAX ALMIN]). APCELL is cell size in φ (degrees), (APCELL ≤ [APMAX APMIN]).
- 4. The minimum and maximum H, λ, φ boundaries of the finite element field, HMIN, ALMIN, APMIN, HMAX, ALMAX, and APMAX. HMIN and HMAX are heights (H) above the reference ellipsoid in meters, (HMIN < 0 and HMAX > HMIN). ALMIN and ALMAX are ellipsoidal longitudes (λ) in degrees, (-180° < ALMIN < ALMAX < 180°). APMIN and APMAX are ellipsoidal latitudes (φ) in degrees, (-90° < APMIN < APMAX < 90°).</p>

These thirteen items are input on four data cards. NORDER, MOBSU, MOBSE, and MOBSN are input in (4I10) format on the first card. The remaining three cards must have a (3E20.14) format, with the finite

element field delimiters, HCELL, ALCELL, and APCELL on the second card, HMIN, ALMIN, and APMIN on the third card, and HMAX, ALMAX, and APMAX on the last card.

Because gravity disturbance observations are generated within PROGRAM LOCALG by SUBROUTINE PTMASS, PROGRAM LOCALG also requires as input the precomputed products of the gravitational constant and each of the mass points of Model 310 and the precomputed rectangular coordinates of each mass point. These quantities should have previously been stored sequentially via an unformatted write on FILE 2. PROGRAM LOCALG accesses these quantities with an unformatted read of FILE 2.

PROGRAM LOCALG produces as output an unformatted sequential file,

FILE 3, which contains the A-matrix for a typical finite element cell

and a set of gravity disturbance observations for each cell of the user

specified finite element field. This file is properly formatted for

later use by PROGRAM FINEG. PROGRAM LOCALG requires two supporting

subroutines - SUBROUTINE CHEBY (or SUBROUTINE ORTHO) and SUBROUTINE

PTMASS.

The basis polynomials are evaluated by the subroutines CHEBY or ORTHO depending on the particular version, and their outputs are used to construct the least squares matrix.

In particular, SUBROUTINE CHEBY (X, N, TA) returns Chebyshev polynomials in array TA, evaluated at a specified value, X, through the specified order, N.  $TA(1) = T_0(X)$ ,  $TA(2) = T_1(X)$ ,  $TA(3) = T_2(X)$ , etc. TA is a seven element array; hence the specified order is bounded,  $0 \le N \le 6$ . SUBROUTINE ORTHO together with SUBROUTINE MULT perform the same function in the case of orthonormal polynomials.

In all versions, the gravity disturbance observations are determined for each cell by SUBROUTINE PTMASS, the final subroutine in Section IIA.

SUBROUTINE PTMASS is used both to produce observations for the least squares fitting process and to produce observations to test against for error analyses of finite element fields. Because SUBROUTINE PTMASS required the precomputed products of each point mass and the gravitational constant and the earth-fixed Cartesian coordinates of each point mass of Model 310, another small program, PROGRAM MASPOS, was necessary to calculate and store these quantities. SUBROUTINE PTMASS uses Mass Model 310 to determine the gravity disturbance observations at each cell grid point.

The rectangular coordinates (x,y,z) of the point are sent to the subroutine and the gravity disturbances are determined by:

$$\delta \underline{\mathbf{g}} = \begin{cases} \delta \mathbf{g}_{\mathbf{x}} \\ \delta \mathbf{g}_{\mathbf{y}} \end{cases} = \sum_{\mathbf{i}=1}^{1080} \frac{PMVALS}{\mathbf{d}_{\mathbf{i}}} \mathbf{i}$$

$$(30)$$

where  ${\tt PMVALS}_{i}$  is the actual value of point mass i which was calculated by MASPOS in Section I

$$\underline{di} = \begin{cases} dx_{i} \\ dy_{i} \\ dz_{i} \end{cases} = \begin{cases} x - POSITS_{i_{1}} \\ y - POSITS_{i_{2}} \\ z - POSITS_{i_{3}} \end{cases}$$
(31)

POSITS are the (x,y,z) coordinates of point mass i which was calculated by MASPOS in Section I and

#### PROGRAM LOCALG LOGIC FLOW (CHEBYSHEV VERSION)

START INITIALIZE CONSTANTS READ POINT MASS VALUES AND FILE 2 PRECOMPUTED POSITIONS INPUT POLYNOMIAL ORDER, NORDER, MOBSU, MOBSE, OBSERVATION GRID PATTERN, MOBSN, HCELL, ALCELL, FINITE ELEMENT CELL SIZE, APCELL, HMIN, ALMIN, AND FINITE ELEMENT FIELD APMIN, HMAX, ALMAX, APMAX BOUNDS GENERATE ALL THE BASIS FUNCTION VALUES AT EACH OBSERVATION GRID POINT WRITE CRITICAL FINITE ELEMENT FIELD INFORMATION FILE 3 ONTO THE SEQUENTIAL ACCESS FILE GENERATE THE LEAST SQUARES MATRIX AND WRITE IT ONTO FILE 3 THE SEQUENTIAL ACCESS FILE LOOP THROUGH THE FINITE ELEMENT FIELD, ONE CELL PER LOOP GENERATE THE GRAVITY ANOMALY OBSERVATIONS (δg<sub>up</sub>,δg<sub>east</sub>,δg<sub>north</sub>) WITH PTMASS WRITE THE THREE OBSER-FILE 3 VATION VECTORS ONTO THE SEQUENTIAL ACCESS FILE GO TO THE NEXT FINITE ELEMENT CELL BY FIRST VARYING φ THEN λ AND FINALLY H UNTIL ALL THE CELLS HAVE BEEN **OBSERVED** 

FIGURE 3

STOP

#### PROGRAM LOCALG LOGIC FLOW (ORTHONORMAL VERSION)

START INITIALIZE CONSTANTS READ POINT MASS VALUES FILE 2 AND PRECOMPUTED POSITIONS INPUT POLYNOMIAL ORDER, NORDER, MOBSU, MOBSE, OBSERVATION GRID PATTERN, MOBSN, HCELL, ALCELL, FINITE ELEMENT CELL SIZE APCELL, HMIN, ALMIN, AND FINITE ELEMENT FIELD APMIN, HMAX, ALMAX, **BOUNDS** APMAX GENERATE THE COEFFICIENTS FOR THE BASIS FUNCTION IN THE H DIRECTION GENERATE ALL THE BASIS FUNCTION VALUES FOR EACH CONSTANT H PLANE GENERATE THE COEFFICIENTS FOR THE BASIC FUNCTION IN THE  $\lambda$  DIRECTION GENERATE ALL THE BASIS FUNCTION VALUES FOR EACH CONSTANT A PLANE GENERATE THE COEFFICIENTS FOR THE BASIS FUNCTION IN THE | DIRECTION GENERATE ALL THE BASIS FUNCTION VALUES FOR EACH CONSTANT & PLANE WRITE THE CRITICAL FINITE ELEMENT FIELD INFORMATION AND FILE 3 ORTHONORMAL COEFFICIENTS ONTO THE SEQUENTIAL ACCESS FILE GENERATE THE LEAST SQUARES MATRIX AND WRITE IT ONTO THE SEQUENTIAL FILE 3 ACCESS FILE

FIGURE 4

## PROGRAM LOCALG LOGIC FLOW (ORTHONORMAL VERSION) CON'T

LOOP THROUGH THE FINITE ELEMENT FIELD, ONE CELL PER LOOP

GENERATE THE GRAVITY ANOMALY OBSERVATIONS (δg<sub>up</sub>, δg<sub>east</sub>, δg<sub>north</sub>) WITH PTMASS

WRITE THE THREE OBSERVATION
VECTORS ONTO THE SEQUENTIAL
ACCESS FILE

FILE 3

GO TO THE NEXT FINITE ELEMENT CELL BY FIRST VARYING
φ THEN λ AND FINALLY H UNTIL
ALL THE CELLS HAVE BEEN
OBSERVED

STOP

FIGURE 4 CON'T

$$d_{i} = dx_{i}^{2} + dy_{i}^{2} + dz_{i}^{2}$$

In particular, SUBROUTINE PTMASS produces gravity disturbance "observations",  $\delta g_{up}$ ,  $\delta g_{east}$ ,  $\delta g_{north}$ , to be used by PROGRAM LOCALG and PROGRAM FINEG for fitting purposes and to be used to test as data against SUBROUTINE FINITE. SUBROUTINE PTMASS produces the gravity disturbance data by using the 1080 point masses of Model 310. SUB-ROUTINE PTMASS requires the COMMON/MASPOS/ be filled with the precomputed products of the gravitational constant and each point mass and the geocentric rectangular coordinates of each point mass. SUBROUTINE PTMASS interfaces with the calling program via COMMON/XYZ/ through which the rectangular coordinates, x, y, z, of the calling point are received and the components of the gravity disturbance  $\delta g_{_{\boldsymbol{X}}}$ ,  $\delta g_{_{\boldsymbol{V}}}$ , and  $\delta g_{_{\boldsymbol{Z}}}$  (DELGX, DELGY, DELGZ) are transmitted. The rectangular components of the gravity disturbance must be transformed to geodetic components  $\delta g_{up}$ ,  $\delta g_{east}$ , and  $\delta g_{north}$ , outside SUBROUTINE PTMASS (i.e., in the calling routine). The logic flow of PROGRAM LOCALG for the different versions are illustrated in Figures 3 and 4.

The Chebyshev and Orthonormal versions both use a second section,

Section IIB to actually calculate the modeling coefficients. These two

versions are different enough to warrant separate discussions.

In the Chebyshev version, the critical field information and the least squares matrix are read in. This version must calculate the modeling coefficients by solving the least squares solution. In lieu of solving the classical normal equations, a more efficient matrix reduction algorithm (ALSQ) is employed. A sequence of Householder elementary matrices are determined which, when premultiplied into the A-matrix

reduce it to a very particular upper triangular form. This triangular system can then be solved by backward substitution for the coefficients.

The A-matrix and its dimensions are passed to the main entry point of SUBROUTINE ALSQ. This section of ALSQ reduces the least squares matrix to its upper triangular form which is returned to FINEG. Next the gravity disturbance component vectors of each cell are read in one at a time. The reduced least squares matrix and the observation vector are then passed to the ALSQ1 entry point of SUBROUTINE ALSQ. This section operates on the observation vectors with the same sequence of Householder transformations, then solves the triangular system for the vectors of coefficients. The process is repeated for each of the up, east, and north measurement data to determine the corresponding coefficient vectors. The coefficients of each cell are written into separate records of a random-access file (FILE 3) to be used by SUBROUTINE FINITE in Section III. In particular, SUBROUTINE ALSQ (A,Y,B,R2,NN,MM,NA) solves the linear least squares problem, ||AB - Y|| = minimum. SUB-ROUTINE ALSQ requires as input the coefficient matrix, A, the observation vector which is to fit, Y, the number of rows used in the A-matrix, NN, the number of columns used in the A-matrix, MM, and the first dimension of the A-matrix, NA. NA must be equal to one plus the maximum number of rows in the A-matrix (NA =  $NN_{max}$  + 1). SUBROUTINE ALSQ returns B, the coefficients of the fit and R2, the sum of the squares of the residuals.

The Orthonormal version of Section IIB produces exactly the same result but the process is much simplier. The simplification is based on the results found in section 5.0 of this report which shows that the

least squares coefficients are given by Eq. (29). Thus, the only processing needed, is to multiply the transpose of the unreduced least squares matrix (i.e., the A-matrix) by the observation vectors (one at a time) to produce the corresponding modeling coefficient vectors (one at a time). Since this was basically the job performed by the secondary entry point, ALSQ1 in the Chebyshev version, the new subroutine for the Orthonormal version is called ALSQ1. (It does not require a secondary entry point.) The coefficients are again returned to the main PROGRAM FINEG and written (in a set of three, one for each component) onto the appropriate record of the random access file (FILE 1). The logic flow charts of both of these versions are given in Figures 5 and 6. In particular, PROGRAM FINEG generates sets of locally valid coefficients Cup, Ceast, and Cnorth for each cell of a finite element field which was previously specified by inputs to PROGRAM LOCALG. The only input required by PROGRAM FINEG is the unformatted sequential file, FILE 3, previously produced by PROGRAM LOCALG. PROGRAM FINEG produces as output a random access file, FILE 1, which contains, in effect, a finite element gravity disturbance field. The first record of the random access file contains certain important information about the finite element field (NORDER, boundaries, etc.). In addition to the first record, the random access file consists of one record of coefficients for each cell of the finite element field. Thus the number of records in the random access file will be equal to the number of cells in the finite element field plus 1. As indicated, PROGRAM FINEG requires one supporting subroutine - SUB-ROUTINE ALSQ.

### PROGRAM FINEG LOGIC FLOW (CHEBYSHEV VERSION)

START

READ CRITICAL FINITE ELEMENT FIELD INFORMATION FROM THE SEQUENTIAL FILE

WRITE CRITICAL FINITE ELE-MENT FIELD INFORMATION ONTO THE FIRST RECORD OF THE RANDOM ACCESS FILE

READ IN THE LEAST SQUARES MATRIX

REDUCE THE LEAST SQUARES MATRIX TO UPPER TRIANGULAR FORM WITH SUBROUTINE ALSO

LOOP THROUGH THE FINITE ELEMENT FIELD ONE CELL PER LOOP

READ GRAVITY ANOMALY
OBSERVATIONS FOR THIS CELL
FROM THE SEQUENTIAL FILE

COMPUTE POLYNOMIAL COEFFI-CIENTS OF THIS FINITE ELE-MENT MODEL WITH ENTRY ALSQ1

WRITE COEFFICIENTS OF THE MODELING EQUATION ONTO THE RANDOM ACCESS FILE

GO TO NEXT CELL

STOP

## PROGRAM FINEG LOGIC FLOW (ORTHONORMAL VERSION)

START

READ CRITICAL FINITE
ELEMENT FIELD INFORMATION
FROM THE SEQUENTIAL FILE

WRITE CRITICAL FINITE ELEMENT FIELD INFORMATION ONTO THE FIRST RECORD OF THE RANDOM ACCESS FILE

READ IN THE LEAST SQUARES MATRIX

LOOP THROUGH THE FINITE ELEMENT FIELD, ONE CELL PER LOOP

READ GRAVITY ANOMALY OB-SERVATIONS FOR THIS CELL FROM THE SEQUENTIAL FILE

COMPUTE POLYNOMIAL COEFFI-CIENTS OF THE FINITE ELEMENT MODEL WITH SUBROUTINE ALSQ1

WRITE COEFFICIENTS OF THE MODELING EQUATION ONTO THE RANDOM ACCESS FILE

GO TO NEXT CELL

STOP

FIGURE 6

Section III, Test Section, is the portion of the software system which evaluates the accuracy and speed of the new modeling approaches verses the Mass Model 310 technique. The main program of this section, FINITES, is the same for both versions of the software which have been developed thus far. The only changes that are necessary are the DIMENSION statements which declare the type and number of the critical field data that must be read from record number 1 (one) of the random access file (FILE 1), the READ statement which performs this operation and the common statements that transfer it to FINITE. Everything else is the same and the procedure is as follows.

After the critical field data are read in, the information produced in Section I for Mass Model 310 is also read in. This information is used to calculate the "actual" gravity disturbance measurements at each test point so that the modeling equation's values at the same points can be compared. Next, the intervals between test points (ISTEPH, ISTEPL, ISTEPP) are read in from cards along with the minimum (HMIN, ALMIN, APMIN) coordinates and maximum coordinates (HMAX, ALMAX, APMAX) of the test region. These coordinates are ellipsoidal. After the validity of the test pattern has been determined, the error analysis is initialized. The ellipsoidal coordinates (H, ALAM, APHI) of each test point are then passed to SUBROUTINE FINITE which returns the three gravity disturbance components which the modeling equation has calculated. These values are saved and the (x,y,z) coordinates of the same test point are determined and sent to SUBROUTINE PTMASS. This subroutine is identical to the subroutine PTMASS discussed earlier. It returns the three gravity disturbance components which Mass Model 310 has calculated. The

disturbance gravity components from PTMASS are first converted to their ellipsoidal values and the components of the two techniques are compared on a point-by-point basis; the error analysis is printed out (the means, standard deviations, and absolute value of maximum errors are printed out along with the average time required to calculate each point).

As noted previously, the actual evaluation of the gravitational disturbances is performed by subroutine PTMASS. For the Chebyshev and Orthonormal versions, although the basis functions are different, the procedure is exactly the same. First the ellipsoidal coordinates (H. ALAM, APHI) of the test point are received from the main program in the common block labelled HLP, and the point is compared against the modeled region minimum (HMIN, ALMIN, APMIN) and maximum (HMAX, ALMAX, APMAX) to see if the test point lies within the region modeled in Section II. If not, the gravity disturbance is evaluated by a call of SUBROUTINE PTMASS, a message indicating this fact is printed and the PTMASS values are returned to FINTES. If the point does lie within the modeled region, the critical field information is used to determine which finite element cell contains the test point. When the cell is determined, the three sets of coefficients (CU,CE,CN) are read from the random access file (FILE 1). Next the coordinates of the test point is normalized with respect to the finite element cell, and these values  $(X_1, X_2, X_3)$  are then passed to SUBROUTINE CHEBY for the Chebyshev version or SUBROUTINE MULT for the Orthonormal version and the basis functions are evaluated at that point. These subroutines are exactly the same as the ones used in Section IIA, to evaluate the basis functions to determine the least squares matrix. The gravity disturbance components are then determined as the sum of the products of the coefficients multiplied by the

evaluated basis function terms. These values are then returned to FINTES for error analysis.

In particular, PROGRAM FINTES is a simple error analysis routine. It compares the gravity disturbance as approximated by SUBROUTINE FINITE with the gravity disturbance as evaluated by SUBROUTINE PTMASS (Model 310). Residuals are printed for each point of a user specified grid, along with the mean of the residuals, the root mean square of the residuals, and the maximum absolute value of the residuals, for each component of the gravity disturbance. The following items are required as input to specify the sample grid:

- The number of samples in each direction, ISTEPH, ISTEPL, and ISTEPP. (ISTEPH > 0, ISTEPL > 0, and ISTEPP > 0). The total number of sample points on the grid will be the product of these three numbers.
- 2. The minimum sample grid coordinate values ("the lower corner"), HMIN (meters), ALMIN (degrees), and APMIN (degrees) and the maximum sample grid coordinate values ("the upper corner"), HMAX (meters), ALMAX (degrees), and APMAX (degrees), (0 < HMIN < HMAX, -180° < ALMIN < ALMAX < 180°, -90° < APMIN < APMAX < 90°).

These nine items are input on three data cards. ISTEPH, ISTEPL, and ISTEPP are input on the first data card in a (315) format. HMIN, ALMIN, and APMIN are input on the second card in (F10.0, 2F10.2) format and HMAX, ALMAX, and APMAX are input on the third card in the same format.

PROGRAM FINTES requires three supporting subroutines - SUBROUTINE FINITE, SUBROUTINE CHEBY or ORTHO (see discussion of these routines above), and SUBROUTINE PTMASS (also discussed above).

Because PROGRAM FINTES uses SUBROUTINE PTMASS, it also requires as input the unformatted sequential file, FILE 2, containing the precomputed products of the gravitational constant and each point mass of Model 310 and the precomputed rectangular coordinates of each point mass.

SUBROUTINE FINITE directly replaces Model 310 in applications involving the gravity disturbance  $\delta g$ . Given the ellipsoidal coordinates, H, ALAM, and APHI  $(H,\lambda,\phi)$ , of a specified point, SUBROUTINE FINITE returns the ellipsoidal components, GU, GE, and GN  $(\delta G_{up}, \delta G_{east}, \delta G_{north})$  of the gravity disturbance approximation,  $\delta G$ . SUBROUTINE FINITE determines which set of finite element coefficients to use at the specified point, reads them into core, and computes  $\delta G$ . SUBROUTINE FINITE requires SUBROUTINE CHEBY or ORTHO (described previously) to produce basis polynomials at the specified point. Whenever the specified point does not lie within the bounds of the finite element field currently accessed, SUBROUTINE FINITE calls SUBROUTINE PTMASS, writes an error message, and sets the error flag ISITIN to zero.

Prior to the first call to SUBROUTINE FINITE, the random access file, FILE 1, must be made available to the calling program, and two flags, IFLAG and ISITIN, must be set to zero and included in a COMMON in both the calling program and SUBROUTINE FINITE. This could be COMMON /HLP/ if necessary (COMMON /HLP/ is the vehicle by which SUBROUTINE FINITE receives the calling coordinates and returns the components of &G), although COMMON /IMARK/ was used.

The following remarks are in order when (1)C FORTRAN is used. The requirements of CDC FORTRAN dictated that the CDC utility subroutines,

subroutines, OPENMS, WRITMS, READMS, and CLOSMS, be used to access the random access file. Furthermore, these CDC utility routines require that the user establish an index array (IMARK, in this case) for the random access file. If an index array is required by a particular version of FORTRAN, the index array, IMARK, should be included in a COMMON in both the calling program and SUBROUTINE FINITE. The logic flow for SUBROUTINE FINITE is displayed in Figure 7.

#### 8.0 Tradeoff Studies

Several tradeoff studies were conducted on the VPI&SU IBM System 360 Model with 158 Processors. The central processor execution times of SUBROUTINE FINITE were compared to SUBROUTINE PTMASS (Model 310). Although it is really a price that is paid only once and a priori in the laboratory, the central processor execution times necessary to create finite element fields were examined. Consideration was given to the tradeoff: total number of finite element coefficients versus NORDER versus cell size, for finite element models of a given region. Lastly, error analyses for a variety of NORDER's and cell sizes were conducted; specifically, the maximum absolute error versus NORDER versus cell size tradeoff was examined.

Table 3 shows the execution time test results for three models:
Chebyshev, Orthonormal, and Pointmass. SUBROUTINE PTMASS has a nearly
constant run time. The run time for SUBROUTINE FINITE, however, depends
on NORDER, the basis polynomial order, and on whether or not a random
access call is needed. A random access call is needed anytime the set
of coefficients currently in core is not the set for the cell in which
the calling coordinates lie; that is, the previous calling coordinates
were in a different finite element cell than the current calling coor-

# SUBROUTINE FINITE LOGIC FLOW

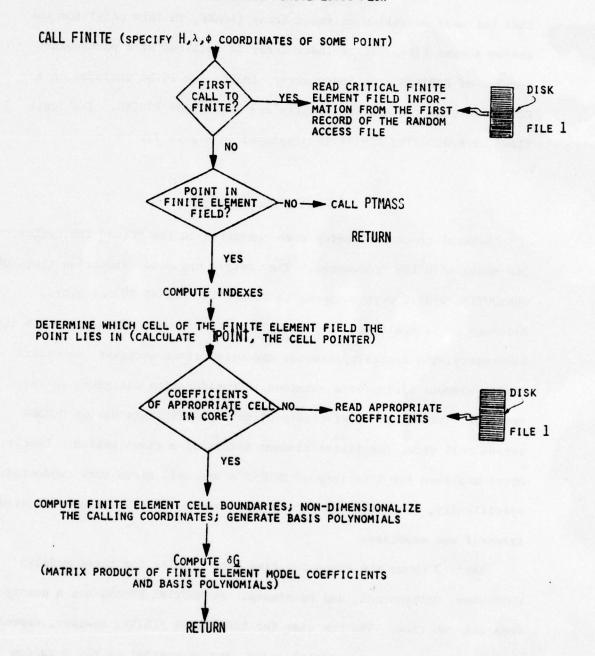


FIGURE 7 SUBROUTINE FINITE LOGIC FLOW

Table 3 Central Processor Execution Time Comparison: Model 310 vs.
Chebyshev Polynomial Model vs.
Orthonormal Polynomial Model (average time per point).

<u>N</u>	Chebyshev	<u>Orthonormal</u>	Pointmass
3	2.2 msec	2.5 msec	230 msec
4	3.0	3.3	230
5	3.9	4.5	230
6	5.4	6.0	230
3	1/105	1/92	Assettic acts 75
4	1/77	1/70	
5	1/59	1/51	
6	1/43	1/38	

dinates. The execution time ratios show that even for the worst cases the Orthonormal model and the Chebyshev model will be at least 40 times faster than an explicit point mass disturbance acceleration model such as Model 310. Note that by using the recurrence relations (24) we can reduce the execution times for Orthonormal so that they are almost identical to the ones produced by Chebyshev.

A certain price must be paid to generate finite element fields. Execution times are shown in Table 4. The price per cell must be multiplied by the total number of cells in the finite element field and added to the price to generate the numbers  $f_{ijk}(\ell)$ . Note that over 90% of the price paid per cell is just the generation of the gravity observations from Model 310. Note the significant time-savings in using the orthonormal model.

The total number of coefficients needed for various finite element field models of the  $35^{\circ} \times 40^{\circ}$  region covered by Model 310 are shown in Table 5. Table 5 assumes that  $H_{cell}$  will be constant at 300 km and that 300 km will be the maximum altitude to be modeled. For a given cell size, it shows how many finite element cells would be needed to model the region and how many total coefficients would be needed for various NORDERS to model the region. Based upon error analyses conducted on a small volume near the center of the  $35^{\circ} \times 40^{\circ}$  region, expected maximum error bounds were determined for certain cell sizes and NORDERS. The upper solid and dashed underline indicates an expected maximum absolute error in mgals. Solid underlines in both cases indicate that the expected maximum absolute error was verified for small test volumes while dashes indicate the apparent trend.

In order to study the maximum absolute error as functions of NORDER and cell size, it was decided to consider the maximum absolute error as

Table 4

Central Processor Times to Create Finite Element Fields

		Cheby	shev		1	Orth	onormal	
N:	3	4	5	6	3	4	5	6
Time to generate matrix A (sec)	.32	.61	1.35	2.90	.41	.70	1.57	3.16
Time to generate the observations using Model	,par,re			SEXES	<sup>9</sup> 80.1			
310 (per cell)	14.87	29.27	50.82	80.12	14.48	28.63	50.43	79.65
Time to process matrix A	1.24	5.45	22.20	77.39	.58	.80	1.29	2.43
Time to calculate the coefficients (per cell)	.43	1.43	3.98	9.23	.22	.75	2.00	5.07
<u>Total Prices</u>				1011				
Time to generate and process matrix A	1.56	6.06	23.55	80.29	.99	1.50	2.86	5.59
Time to generate obser- vations and coefficients				144	97.2			
using Model 310 (per cell)	15.30	30.70	54.80	89.35	14.70	29.38	52.43	84.72

Table 5 Total number of Coefficients vs. N. vs. Cell Size For a 300 km  $\times$  35 $^{\circ}$   $\times$  40 $^{\circ}$  Finite Element Field. (Chebyshev and Orthonormal)

Cell Size $(\lambda_{\text{cell}} \times \phi_{\text{cell}})$	No. of Cells	3	4	5	6
1.0° × 1.0°	35×40 = 1400	84,000	147,000	235,200	352,800
1.25° × 1.25°	28×32 = 896	53,760	94,080	150,528	225,792
1.5° × 1.5°	24×27 = 648	38,880	68,040	108,864	163,296*
1.75° × 1.75°	20×23 = 460	27,600	48,300	77,280	115,920
2.0° × 2.0°	18×20 = 360	21,600	37,800	60,480	90,720
2.25° × 2.25°	16×18 = 288	17,280	30,240	48,384	72,576
2.5° × 2.5°	14×16 = 224	13,440	23,520	37,632	56,448
$2.75^{\circ} \times 2.75^{\circ}$	13×15 = 195	11,700	20,475	32,760	49,140
3.0° × 3.0°	12×14 = 168	10,080	17,640	28,224	42,336

Given:  $H_{Cell} = constant = 300 \text{ km}$ 

<sup>\* =</sup> Finite element fields above this line are expected to have  $\left| \text{error} \right|_{\text{max}} \leq 1.5 \text{ mgals.}$ 

<sup>+</sup> = Finite element fields above this line are expected to have  $\left|\text{error}\right|_{\text{max}} \leq 3.0 \text{ mgals.}$ 

a function of cell size only for fixed NORDER (see Figure 8) and to consider maximum absolute error as a function of NORDER for fixed cell sizes (see Figure 9). For all cases, cell altitude ( $H_{cell}$ ) was fixed at 300 km.

Figure 8 shows that the maximum absolute error varies approximately as the square of the cell dimension in  $\lambda$  or  $\phi$ . For constant cell altitude and for small  $\lambda_{\text{cell}}$  and  $\phi$ , the volume will vary approximately as the square of  $\lambda_{\text{cell}}$  of  $\phi_{\text{cell}}$ , if  $\lambda_{\text{cell}} = \phi_{\text{cell}}$ . Hence the maximum absolute error is really proportional to cell volume. Figure 9 shows that the maximum absolute error varies approximately inversely as the square of NORDER. As the number of coefficients per cell varies approximately as the square of NORDER, the maximum absolute error may be considered to vary nearly inversely with the number of coefficients per cell.

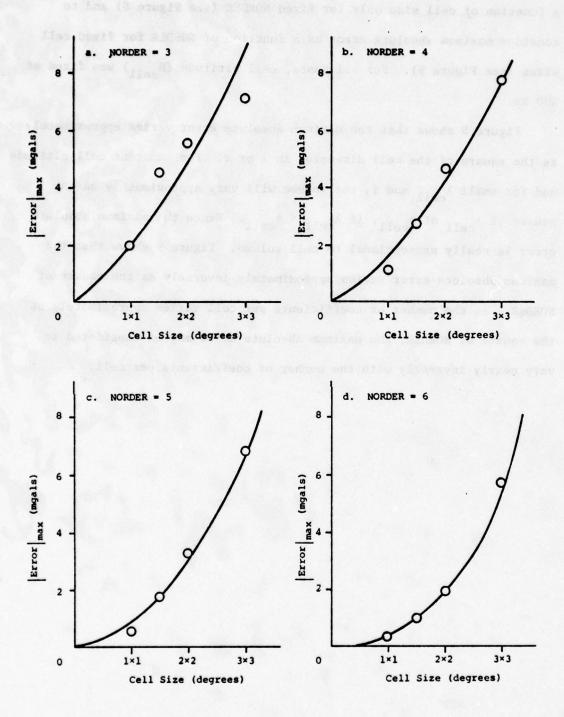


Figure 8 Maximum Absolute Error versus Cell Size for Fixed NORDER's.

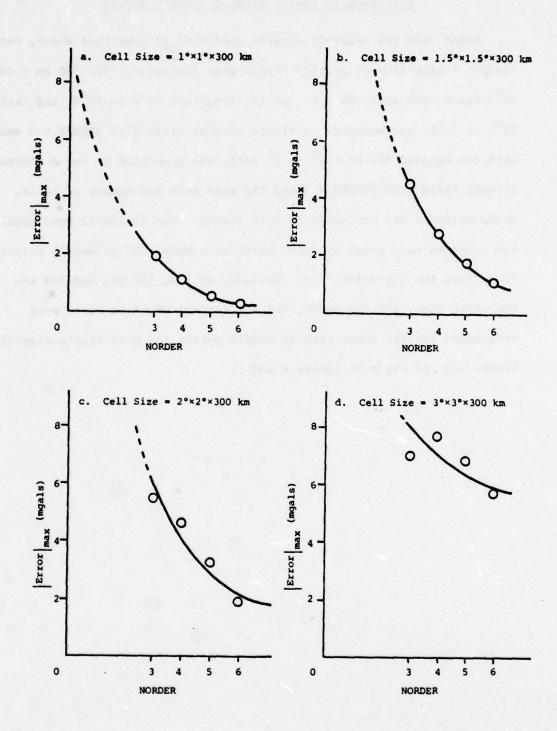


Figure 9 Maximum Absolute Error versus NORDER for Fixed Cell Sizes.

## 9.0 Example Finite Element Gravity Models

Based upon the tradeoff studies conducted as described above, two example finite element gravity fields were generated. The 300 km  $\times$   $10^{\circ}$   $\times$   $10^{\circ}$  region with altitude 0 to 300 km, longitude  $70^{\circ}$ E to  $80^{\circ}$ E, and latitude  $25^{\circ}$ S to  $35^{\circ}$ S was modeled. A finite element field with NORDER = 3 and with one hundred 300 km  $\times$   $1^{\circ}$   $\times$   $1^{\circ}$  cells was generated as was a finite element field with NORDER = 5 and the same size and number of cells. An error analysis was conducted on both fields. The finite element model was compared on a point by point basis with Model 310 at sample points "away from the fit points" for altitudes of 1 m, 150 km, and 299 km. The error mean, the RMS error, and the maximum absolute error were determined for the three sets of sample points for both finite element fields and are shown in Tables 6 and 7.

Table 6

N=3 Example Finite Element Gravity Disturbance Field (Chebyshev and Orthonormal)

	δg <sub>up</sub> (mgals)	<sup>δg</sup> east (mgals)	δg <sub>north</sub> (mgals)
H = 1 m			
ERROR	005	001	.004
RMS ERROR	.963	.645	.727
ERROR  <sub>MAX</sub>	2.717	2.074	2.236
<u>H = 150 km</u>			
ERROR	023	.003	.009
RMS ERROR	.352	.224	.270
ERROR   MAX	1.149	.701	1.039
H = 299 km			
ERROR	.018	006	007
RMS ERROR	.424	.277	.322
ERROR  <sub>MAX</sub>	1.130	.771	1.002

N=5 Example Finite Element Gravity Disturbance Field (Chebyshev and Orthonormal)

	δg <sub>up</sub> (mgals)	<sup>δg</sup> east (mgals)	δg <sub>north</sub> (mgals)
H = 1 m			
ERROR	004	.001	.002
RMS ERROR	.224	.150	.167
ERROR  <sub>MAX</sub>	.918	.637	.581
H = 150 km			
ERROR	.007	002	004
RMS ERROR	.071	.044	.055
ERROR  <sub>MAX</sub>	.216	.157	.215
H = 299  km			
ERROR	002	.001	.001
RMS ERROR	.101	.067	.077
ERROR  <sub>MAX</sub>	.329	.254	.249

## 10.0 Orthogonal Approach

The main purpose of this section is to indicate a way to produce an accurate model for the geopotential itself. The Chebyshev and Orthonormal approaches model the gravity disturbance acceleration components independently and do not take into account the fundamental functional relationship

$$g = \nabla U \tag{32}$$

Therefore, no rigorous representation for U can be obtained by using gravity components derived either by the Chebyshev or Orthonormal approach.

A finite element procedure for modeling the geopotential that takes the functional relationships (32) between the geopotential and its gradient into account appears desirable. Not only will this enable us to obtain an approximation for U, but a significant reduction in the overall cost of determining gradient models is anticipated.

Let us consider the following preliminary approximation of the geopotential disturbance function

$$\delta U = \sum_{n=0}^{N} \sum_{i=0}^{n} \sum_{j=0}^{i-n} c_{ijk} f_{ijk} , k = n-i-j$$
 (33)

where  $c_{ijk}$  are modeling constants to be determined, and  $f_{ijk}(x,y,z)$  are arbitrary basis functions. Using Eq. (32) and (33) we can write the acceleration disturbances as follows

$$\delta g_{x} = \sum_{n=0}^{N} \sum_{i=0}^{n} \sum_{j=0}^{i-n} c_{ijk} f_{ijk}$$
(34a)

$$\delta g_{y} = \sum_{n=0}^{N} \sum_{i=0}^{n} \sum_{j=0}^{i-n} c_{ijk} f_{ijk}$$
(34b)

$$\delta g_{z} = \sum_{n=0}^{N} \sum_{i=0}^{n} \sum_{j=0}^{i-n} c_{ijk} f_{ijk}$$
(34c)

with

$$f_{ijk} = \frac{\partial}{\partial x} f_{ijk}, f_{ijk} = \frac{\partial}{\partial y} f_{ijk}, f_{ijk} = \frac{\partial}{\partial z} f_{ijk}$$
 (35)

The local modeling coefficients  $c_{ijk}$  are to be determined via a least squares approximation technique, taking into account the local observations of all three gravity components  $(\delta g_x, \delta g_y, \delta g_z)$  simultaneously.

The normal equations can be written in the usual manner

$$\underline{\mathbf{c}} = (\mathbf{A}^{\mathrm{T}}\mathbf{A})^{-1} \mathbf{A}^{\mathrm{T}} \delta \underline{\mathbf{g}}^{*} \tag{36}$$

where

$$\underline{\mathbf{c}} = \begin{cases} \mathbf{c}_{000} \\ \mathbf{c}_{001} \\ \mathbf{c}_{010} \\ \cdots \\ \mathbf{c}_{N00} \end{cases}, \quad \delta \underline{\mathbf{g}}^* = \begin{cases} \delta \mathbf{g}_{\mathbf{x}}^*(1) \\ \delta \mathbf{g}_{\mathbf{y}}^*(1) \\ \delta \mathbf{g}_{\mathbf{z}}^*(1) \\ \cdots \\ \delta \mathbf{g}_{\mathbf{x}}^*(\mathbf{m}) \\ \delta \mathbf{g}_{\mathbf{y}}^*(\mathbf{m}) \\ \delta \mathbf{g}_{\mathbf{z}}^*(\mathbf{m}) \end{cases}$$

$$(37)$$

are the column matrix of modeling constants c ijk and the column matrix of gravity disturbance observations, respectively. Furthermore,

in which m indicates the total number of local measurements of the disturbance acceleration.

It is now possible to evaluate the constants  $c_{ijk}$  from Eq. (36) in the same manner as the Chebyshev approach. However, we wish to introduce a significant simplification by using orthogonal basis functions, similar to the Orthonormal approach. To this end, let us write the expression for a generic element of the matrix  $A^T\!A$  as follows

$$\sum_{\ell=1}^{m} F(\ell)$$

$$\ell=1 \text{ ijk}, \alpha\beta\gamma$$
(39)

where

$$F(\ell) = f_{ijk}(\ell) f_{\alpha\beta\gamma}(\ell) + f_{ijk}(\ell) f_{\alpha\beta\gamma}(\ell) + f_{ijk}(\ell) f_{\alpha\beta\gamma}(\ell) + (40)$$

Let us now introduce basis functions of the form

$$f_{ijk}(x,y,z) = f_i(x) f_j(y) f_k(z)$$
 (41)

so that,

$$\mathbf{x} \\
\mathbf{f}_{\mathbf{i}\mathbf{j}\mathbf{k}} = \mathbf{f}_{\mathbf{i}}^{\prime}\mathbf{f}_{\mathbf{j}}\mathbf{f}_{\mathbf{k}} , \quad \mathbf{f}_{\mathbf{i}\mathbf{j}\mathbf{k}} = \mathbf{f}_{\mathbf{i}}\mathbf{f}_{\mathbf{j}}^{\prime}\mathbf{f}_{\mathbf{k}} , \quad \mathbf{f}_{\mathbf{i}\mathbf{j}\mathbf{k}} = \mathbf{f}_{\mathbf{i}}\mathbf{f}_{\mathbf{j}}\mathbf{f}_{\mathbf{k}}^{\prime}$$
(42)

where the derivatives denoted by primes are taken with respect to the argument. It is clear that the generic element (39) can be written as

$$\sum_{\mathbf{x}} \mathbf{f}_{\mathbf{i}} \mathbf{f}_{\alpha} \sum_{\mathbf{y}} \mathbf{f}_{\beta} \sum_{\mathbf{z}} \mathbf{f}_{\mathbf{k}} \mathbf{f}_{\alpha}$$

$$+ \sum_{\mathbf{x}} \mathbf{f}_{\mathbf{i}} \mathbf{f}_{\alpha} \sum_{\mathbf{y}} \mathbf{f}_{\beta} \sum_{\mathbf{z}} \mathbf{f}_{\mathbf{k}} \mathbf{f}_{\alpha}$$

$$+ \sum_{\mathbf{x}} \mathbf{f}_{\mathbf{i}} \mathbf{f}_{\alpha} \sum_{\mathbf{y}} \mathbf{f}_{\beta} \sum_{\mathbf{z}} \mathbf{f}_{\mathbf{k}} \mathbf{f}_{\alpha}$$

$$(43)$$

where the respective summations run over all the observations in the respective directions.

Next, let us assume that the sets of functions

$$\{f_{i}\}$$
 ,  $\{f_{i}\}$  ,  $\{f_{k}\}$ 

are orthonormal over the interval of validity and that similarly

$$\{f_{i}^{!}\}$$
 ,  $\{f_{i}^{!}\}$  ,  $\{f_{k}^{!}\}$ 

are orthogonal with respect to the same interval, then it is clear that  $\mathbf{A}^T\mathbf{A} = \mathbf{D}$  becomes diagonal with diagonal elements  $\mathbf{D}_{ijk}$  given by

$$D_{ijk} = \sum_{x} f_i^{2} + \sum_{y} f_j^{2} + \sum_{z} f_k^{2}$$

Equation (36) becomes

$$c = D^{-1}A^{T} \delta g^{*}$$

or, more explicitly,

$$c_{ijk} = D_{ijk}^{-1} \sum_{\ell=1}^{m} (f_{ijk}^{x} \delta g_{x}^{*} + f_{ijk}^{y} \delta g_{y}^{*} + f_{ijk}^{z} \delta g_{z}^{*})_{\ell}$$
(44)

The only remaining task is to find an orthonormal set of functions  $\{f_i(x)\}$  such that the derivative set  $\{f_i(x)\}$  is also orthogonal over the

same interval of validity. Note that the same set can be used in all three directions x,y,z.

To date, only NORDER = 3 has been tried. As expected, however, not enough accuracy is obtained and further work is underway for higher NORDERs. It is expected that the number of modeling constants  $c_{ijk}$  will be significantly reduced compared to the total number of modeling constants  $c_{ijk}^{\ \phi}$ ,  $c_{ijk}^{\ \lambda}$ ,  $c_{ijk}^{\ \gamma}$ , necessary to model the gravity disturbance components independently. This fact, not only increases the computational speed of the algorithm, but also reduces the storage requirements. Moreover, it is now possible to obtain a valid approximation for the geopotential itself.

## 11.0 References

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<sup>\*</sup>These Orthonormal programs and subroutines represent minor modifications of the corresponding Chebyshev Polynomial Software.

<sup>\*\*</sup>The numerical results (for fixed NORDERS) are identical for the two approaches. The orthonormal vs. Chebyshev tradeoffs are with regard to efficiency, not accuracy, see the above discussion in Section 8.0.

MODEL 310

POINT MASS DISTURBANCE ACCELERATION CALCULATION

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                                                                                                                                                                                                                       RNI=AZ/USEKI(AZ#LUSP#CUSF + BZ#SINF)
TEMP=(RNI+RHI)#CGSP
TEMPZ#BZ#RNI/AZ+RHI
ALAMEM=3575.00
DU ZO J=1*36
                                                                                   FISTITION OF THE 1080 MASSES
                                                                                                                                                                                                                                                                                                                                                                            PMVALS(1PT)=KM1*8166*1.019
                                                                                                                                                                                                                                                                                                                                                                                                                 POSITS(IPT, 1) = TEMP*COSL
PCSITS(IPT, 2) = TEMP*SINC
PCSITS(IPT, 3) = LEMPZ*SINP
CONTINUE
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                   WKITE(2) PMVALS, PUSITS
                                                        FACM A STARTING PUTNI UF:
                                                           35750
85750
8755
                                                                                                                                                                                                                                                            DC 20 J=1,36
ALAMEM=ALAMEM+50.DC
ALAME=ALAMEM/KADMIN
SINL=DSIN(ALAME)
CCSL=UCUS(ALAME)
MAIN
                  READ IN THE MASS CODES
                                     READ (5, 403) MASSES
                                                                                                                                                     PHIME—875.00
20 20 1=1.50
PHIMEPHIM-50.00
PHIEPHIM/RADMIN
SINY=USIN(FHI)
ULSY=UCUN(FHI)
                                                                                                                                                                                                                                                                                                                     1P1=1P1+1
M1=MASSES(1P1)-2
RM1=M1
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                            1=1,20
                                                                                                                 A.=A*A
32=E*D
Khl=-Eucuu.Du
                                                                                      CALCULATE THE
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0+0300000000	7046148100140	0+05まんで、82717	0+0/12/0/0/0/0/0/0/0/0/0/0/0/0/0/0/0/0/0/0
.e700000000+0	. ce84404356D+0	-2574438707D+U	Z-18308498470+C
٠	-+116423534D+C	0+0450205086447	Z-18508498470+C
0+00000000 E9.	0+34161460+00	. 50341224550+0	0+118308478470+C
- a 100000000000000000000000000000000000	•4565551553U+0	36470807340+0	2 - 18308478470+0
0+0000000000000000000000000000000000000	3763364560+0	0+00/1000000000000000000000000000000000	4 - 183084784 10+0
	04090466666	-4588846551U+0	0+07+0440401-7
2+03 30000012 4 4	0.0000000000000000000000000000000000000	- 4.0.2.0.0.0.0.0.0.0.0.0.0.0.0.0.0.0.0.0.	こうことものののはいい
0+0.00000000000000000000000000000000000	2+174 × (1-174)	のでは、これのでは、これでは、これでは、これでは、これでは、これでは、これでは、これでは、これ	の中心でもよりないであれてい
0+0000000000	ロキロのなくのないとかった。	0+07/209/204/04	の中のできないののできる。
04 000 000 000 L9	67608374450+0	07101454920+0	0-15308478470+0
-67000000000000-c	-8105577495U+O	.618588675eD+0	2-18308498470+0
0	.73465143720+0	.04441426730+0	2.18308458470+0
0+0000000019.	.c5237c18300+0	0+051985+0599	2-18308478470+0
פיפונת ממם בהם +ם	340916516160+6	0+05/472/4759.	2-16208478470+0
1+00000000000	1+01T+0+A100+•	114 130 40 80 040	0+01+01+01-1 0+01-1
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0+000000000000	0+126255380+0	0+05557054997	2-26822723450+0
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CHEBYSHEV POLYNOMIAL FINITE ELEMENT SOFTWARE

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WKITE(3) HMIN, HMAX, HCELL, ALMIN, ALMAX, ALCELL, APMIN, APMAX, APCELL, NH, +NLAM, NPHI, NCKDEK, NC, MUDS, NCELLS
                                                                         WRITE(6,4) NUKDER,NC
WRITE(6,4) MUBSU,MUBSE,MUBSN,MOBS
WRITE(6,5) NUELL,ALCELL,APCELL,HMIN,ALMIN,APMIN,HMAX,ALMAX,APMAX
      20/33/12
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                 EVALUALE INC. CASIS FUNCTION AT EACH GRID PUINT FUR EACH UNDER OF FUNCTIONS OF TO THE NECESSARY URDER
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                CLUMPULE THE NUMBER OF FINAL CELLS IN UP-EASTERN, AND NORTHERN DIRECTION AND THE TOTAL NUMBER OF FINAL CELLS IN THE FINITE LLEMENT FIELD.
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                             DETERMINE THE NORMALIZED LAMBA CODDINATE FOR EACH GRID POINT
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                            DEJERMINE THE NURMALIZED IN COURDINATE FUR EACH GKID FUINT
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                               WKITE FILLD DATA UNTO FIRST RECORD OF SEQUENTIAL FILE
          DATE = 79206
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                COMPUTE THE INCREMENT SIZE MITHIN EACH CELL.
                                                                                                                                                                                                                                      LUNVER! ANGLES FROM DECREES TO RADIANS
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                          KNN= IMMAX-HMIN J/HCELL+.9990G
KNL= (ALMAX-HALMIN J/ALLELL+.9990G
KNN= (AFMAX-AFMIN J/AFCELL+.9990G
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                            ALAM=ALMIN-DLAM
DL 150 1L=1,MubSe
ALAM=ALAM+ULAM
XZ=(ALAM-ALMIN)/ALCELL
CALL CHESY(X2,NUNDEK;[Y(1,1L))
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                   XI=(n-nMIN)/nctll
CALL CHEBY(XI)NUKLEK, IX(I,1H))
CGN/INUE
                                                                                                                                                                                                                                                                                                     ALCELL=ALCELL+DEGKAU
APCELL=ACELL+DEGKAU
ALMIN=ALMIN*DEGKAU
APMIN=AFMIN*DEGKAU
ALMAX=ALMAX*DEGKAU
AFMAX=ALMAX*DEGKAU
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NLIE: IF A SYMMEIKIL GRID PAITENN IS ALWAYS CHUSEN ISUCH AS +*+*+)
THE INNEE SETS OF EVALUATIONS ABOVE ARE REDUNDANT AND TWO CLULD BE ELIMINATED
   20/33/12
                                                                                                                                                               DETERMINE INC NURMALIZED PHI CLURLINATE FUR EACH GRID POINT
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                              KESET THE LAMBA CCURUINATE TO 113 CELL MINIMUM VALUE
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                INCKEMENT THE LAMBA CUBNIDATE TALMUEL! I LELL SIZE
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   DATE = 79206
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                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                  A (MJ, 11)=TX(NX, 1H)*IY(NY, 1L)*1L(NZ, 1P)
CEN LAJE
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                    WRITE THE BASIS FUNCTION MAIKIX 10 DISK
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                                                                                                                                                                                                                                                                                                                                                   X3=(AFHI-AFMIN)/AFCELL
CALL CHEBY(X3,NORDER, 12(1,1P))
CON 11NUE
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                GENERATE DATA FUR LACH CELL.
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                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                  UC 150 IN=1,MC65C
UC 150 IC=1,MC65C
UC 150 IY=1,MC65N
AL=MJ+1
                                                                                                                                                                                                                                           APPLIMATMIN-UPPLI
UC 140 IPPLIMEDSN
APPLIMATHI-UPPLI
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MAIKIX FOR 3-2 KUIAIIUN (LAMBDA,-PHI) FRUM X,Y,Z, TU
 20/33/12
                                                                                                                                                                                         INCREMENT THE LAMBA CCORDINGTE (ALMCEL) I GRID INCREMENT SIZE ALAM=ALAM+ULAM
CCSC=CCS(ALAM)
SINC=USIN(ALAM)
AFRI=AFMCEC-UFFIL
UC 250 IP=1,MCBSN
                                                                                                                                                                                                                                                                                                                                  TRANSFORM ALL ELLIPSUIDAL COURDINATES INIU X,Y, 2 CCURDINATES
                                                                                                                                                                                                                                                              INCREMENT THE PHI CUURLINATE LAPMERT I GRID INCREMENT SIZE
                                                                                                                           INCREMENT THE H CUURDINATE IMMUELL) I GRID INCREMENT SIZE n=H+DH
                                                                                                                                                                                                                                                                                                                                                                                                                                             UBIAIN THE GRAVITY "LESCHVALIEN" AT IMIS GRID PUINT
DATE = 79206
                                                            INCREMENT THE PHT COURDINATE LAPMCEL! I LELL SILE
                                                                                                                                                                                                                                                                                                                                                      + BZ#SINP#SINP
                                                                                                                                                                                                                                                                                                                                                      RN=DSERTIAZ*LOSP*CUSF
 MAIN
                                                                               APMLEL=APMLEL+APCELL
                              APMCEL=APMIN-APCELL
UU 300 IPC=1,NPH1
                                                                                                HEHMCELL-DH
DE 250 IH=1,MUBSU
                                                                                                                                                   ALAM=ALMCEL-ULAM
                                                                                                                                                                                                                                                                                  APRI PAPRI + UPRI
COSP = UCCOS ( PPRI )
SINY = USIN( APRI)
RU= MU+1
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                     UIKECTION CUSINE
UP, EASI, NURIH
                                                                                                                                                                                                                                                                                                                                                                           T=(KN+H)*CUSF
                                                                                                                                                                                                                                                                                                                                                                                                                 13154 (H+27) =
                                                                                                                                                                                                                                                                                                                                                                                                                                                                   CALL FIMASS
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                               C11=C0SP*C0SL
C1Z=C0SP*51NL
C13=S1NP
C21=-51NL
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RELEASE 2.0
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FURTRAN IV G1
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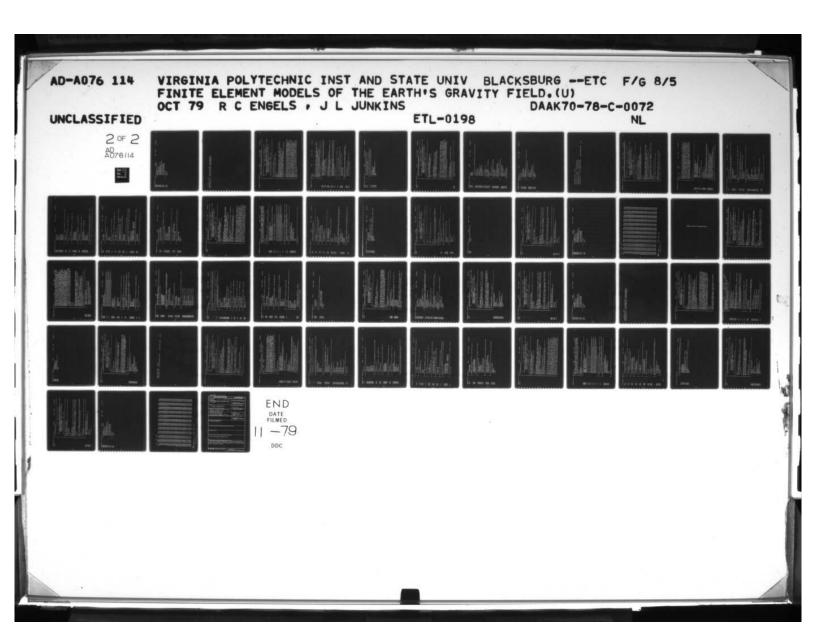
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                                                                                                                                                                           WAITE THE THREE UBSERVATION ARRAYS UNTO THE SECUENTIAL FILE
DATE = 79206
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                                                                                                    SICKE THE UDSERVATION DISTURBANCES
                                                                                                                                                                                      WAITE(3) (DELGU(1),1=1,Mubs)
waite(3) (UELGE(1),1=1,MUbs)
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                                                                                                                       ACO DELGO(MJ)=GU
DELGE(MJ)=GE
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ZEC CUN)INOE
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C32=-51NP*51NL
C53=CC5P
KELEASE 2.0
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FURTRAN IV G1
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ORTRAN IV GI . RE	RELEASE 2.0 CHEBY DATE = 79206 20/33/12
1000	SUBROUTINE CHEBY(Xon+TA)
	C C C DEPARIMENT OF JUNKINSTREMI C. ENGELS, AND JOHN J. SMITH C SCHOOL OF FOLTERING SCIENCE AND MECHANICS SCHOOL OF ENGINEERING SCIENCE AND MECHANICS VIKEINIA FOLTIECHNIC INSTITUTE AND STATE UNIVERSITY C BLACKSCOKG, VIKEINIA 24001
	UATE
	INPUIS
	C FROM SUERGUTING CALL. C X-THE NUMMALIZED CUCRLINATE WHERE THE FUNCTION IS IN SE EVALUATED. C N-Chiem of the SASIS FUNCTION JESTRED
	· Pruess
	C LECKLINALE VALUE AND AND THE BASIS FUNCTION AT THE NUMBELIZED (N). C LECKLINALE VALUE X. TUR EACH URDER UF THE FUNCTION SPECIFIED (N). THE FUNCTION VALUES ARE RETURNED IN VECTOR IA.
	בוסאומו אינור אינ
	C IL SUBREULINE RETURN. C IA-TUNCTIUN VALUES. CF. URDER N EVALUATED AT X
2003	JMPL1011 REAL*8 (A-F,U-Z)
4 c o c c c c c c c c c c c c c c c c c	JE (N.6). C. AND. N. Ll. 7) GUIU 10 WELLE (C.) N PURMAI(Z4MUILLE GAL NURUEK: 10) STUF 10 LUNIANDE
	C TRANSFORM THE CUCRDINALE FANCE (0,1) IC (-1,1)
\$0000000 \$0000000000000000000000000000	C XEAR=2.50*X-1.00 [A(1)=1.50* [A(2)=XBAR NP1=N+1 UG 20 1=3,NP1 20 1A(1)=2.50*XBAN*1A(1-1)-1A(1-2) REIURN

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20/33/12					
DATE = 7920c					
RELEASE 2.0 FIMASS	10 CCN 11NGE UX=PCS175(11) -X	U2=FUSITS(I1:5)-Z U1S1SG=UX*UX+UY+UZ*UZ D1ST=DSQKI(U1S1SQ) TEMP=PMVALS(1)/U1S1/U1S1SQ	DELGX=DELGX+(DX*TEMP) JELGY=DELGY+(DY*1cMP) DELGZ=DELGZ+(DZ*1cMP)	C CONTINUE	RETURN
FORTRAN 1V G1	5000	17500	0017 5018 5019	0200	7790 0057

MUDIFICATIONS BY JUHN L. JUNKINS, KEMI C. ENGELS, AND JUHN J. SMITH DEPARIMENT OF ENGINEERING SCIENCE AND MECHANICS SCHOOL OF ENGINEERING SCIENCE AND MECHANICS SCHOOL OF ENGINEERING VIRGINIA POLYTECHNIC INSTITUTE AND STATE UNIVERSITY ELACKSBORG, VIRGINIA 24561 THIS PRECEASE ACCEPTS FROM DISK, INPUT DESCRIBING THE LIMITS OF THE REGION TO BE MUDELED, THE CRUEK OF THE EASIS FUNCTIONS TO DE USED IN THE MODELING, A LEAST SQUAKES MATRIX MADE UP OF DASIS FUNCTION PRODUCED AND SETS OF THREE VECTORS OF GRAVITY ANDMALY OBSERVATIONS FOR THE FACT. CELL TIME NOT THE NEUTON OF DASIS FUNCTIONS AND THE FIRST THE LOSS FILE TO THE LEAST SQUAKES MATRIX IS THE NEUTON OF SANDOM ACCESS FILE AND THE AND THE SECOND SECTION OF ALSO THE THE NEUTON OF ALSO THE DASIS OF THE OBSERVATION OF ALSO THE BASIS OF CELL. ARE THE UBSERVATION OF ALSO THE BASIS OF CELL. ARE THEN WAITHEN INTO THE NEUTON OF ALSO THE BASIS AND CELL ARE THEN WAITHEN INTO THE NEXT RECEND SETS THE SANDOM ACCESS THE SANDOM ALCOS TILE. SINCE THE CURTICLEN'S THE NEXT RECENDED SET THE THE NEXT TO THE THE NEXT TO THE NEXT TO THE NEXT TO THE NEXT TO THE THE NEXT TO THE TOWN TO THE NEXT TO THE 20/33/39 FREM DISK FILE 3 HMIN, ALMIN, AFMIN-MINIMUM BUUNDS UP FINITE ELMENT FIELD DESIRED HMAX, ALMAX, AFMEAL-MAXIMUM BUUNDS UF FINITE ELEMENT FIELD DESIRED HNEAL, ALCELL, FREEL-LEL SIZE IN H, LAMBUA, PHI NCHLE, ALCELL, FREEL-LELS IN H, LAMBA, PHI DIRECTION NCRDER-CRDER LF PULYNOMIALS UPSIRED NCRDER-CRDER LF PULYNOMIALS UPSIRED NC-NUMBER UF CUEFFICIENTS IN THE MUDELING EQUATION NC-NUMBER UF CUEFFICIENTS IN THE MUDELING EQUATION AL-DASIS FUNCTION MAIRIX FOR LEAST SQUARES FIT DELGU, PELLE, DELGN-CRAVITY UISTURGANGE GUMPUNENTS IN H, LAMBA, PHI SECTION 113 SUBER -- CONFIDENT DETERMINATION PROCKEM JUHN L. JUNKINS AND JUHN SAUNDERS UATE = 79206 PROCKAM FINEGLOUIPUI, TAPED=OUIPUT, LAPELI, TAPES) FLR F. SUME FLINI IN SPACE A.Y.Z.—L.ARIM-FIXED AFLIANGULAR CUCKLINAIES M.FLAM, FFLITT, LLIFSCIERE CUCKCINAIES NORDERT-UNDER IF INE DASIS FUNCIIUNS DESIRED 1,1579 UF LAST MUDIFICATION -- APRIL MALR ANCHALY COMPOSENIO. トドしてにひい DATE RELEASE 2.0 2 FURTRAN IV GI

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CALL TIMECN
READIS) HMIN, HMAX, HCELL, ALMIN, ALMAX, ALCELL, APMIN, APMAX, APCELL, Wh,
*NLAM, NPHI, NUH UER, NC, MUDS, NUELLS
WKI I E (1 * 1) HMIN, HMAK, HCELL, ALMIN, ALMAX, ALCELL,
*NRI E (1 * 1) HMIN, HMAK, HCELL, ALMIN, ALMAX, ALCELL,
*NH, NLAM, NPHI, NUKDLK, NC
                                                                                                                                                                                                                                                                                                                                                                                              IMPLICIT REAL#G (A-H,C-2)
COMMUN /FLUATA/HMIN, HMAX, HCELL, ALMIN, ALMAX, ALCELL, APMIN, APMAX,
*APCELL, NH, NLAM, NPH1, NUCKUE K, NC
DIMFNSION A (344,85), DELG(343)
DIMFNSION A (344,85), DELG(343)
DIMFNSION COLONI, CE (64), CN(64)
CATE MAXUEL (544,65)
                                                                                                                                          HMIN-ALMIN-FPMINT-MINIMUM OLUNDS OF FINITE ELMENT FIELD DESIRED MAX, ALMAX, APMAX—MAXIMUM GUUNDS OF FINITE ELEMENT FIELD DESIRED HOLELL, ALCELL, APCELL—TCELL SIZE IN M. LAMBDA, PHI NELD DESIRED NH, NPHI—NOMBER OF CELLS IN M. LAMBA, PHI DIRECTION NC—NOMBER OF COLFTICIENTS IN THE MODELING EQUATION COMPONER OF COLFFICIENTS IN THE MODELING EQUATION OLD COMPONENTS OF CULCEFFICIENTS FOR THE M. LAMBDA, PHI COMPONENTS OF THE CRAVITY ANGMALY
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                         FUR EACH CELL, READ UNE SET UF GRAVITY UBSERVATIONS AND CUMPULE THE CUEFFICIENT AKKAY FUR EACH CUMPUNENT UF GRAVITY
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                             REDUCE THE LEAST SAUAKES MATRIX TO UPPER INJANGULAR FURM
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READIS) (LELE(1),1=1,MCDS)
CALL ALSU1(A, LeLG, CU, SCMS2, MOBS,NC, MAXUEl)
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                             CALL ALSGIA, DELG, CU, SUMSQ, MUBS, NC, MAXOBI)
CALL TIMECK (ICP)
FIME=ICP/ICC, DC
WKITE(0, 7) FIME
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CALL ALSUITA, DELG, CE, SUMSG, MOBS, NC, MAXUB1)  CEAL ALSUITA, DELG, CH, SUMSG, MUES, NC, MAXUEL)  CALL ALSCITA, DELG, Ch, SUMSG, NC, MAXUEL)  COZT  CALL CONTINUE  COZZ  CAL CONTINUE  CALL TIME ETCP I CUITC), CHITC), CHITC), CHITC)  TIME ETCP I CONTINUE  COZZ  CALL TIME CKITCP)  TIME ETCP I CONTINUE  TANDER CKITCP)  TANDER CKITCP  TA	GRTRAN IV GI KELEASE 2.0	NICH	UATE = 79206	20/33/3
<b>.</b>	CALL ALSU	II A, DELG, CE, SUMSQ, MO	BS,NC, MAXUBI)	
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ר ר ז	L STURE EACH LE	DEFFICIENT ANNAY IN	A KANDUM ALLESS FILE	
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UMPLICIT REALTS (A-H)-L-1)
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10 SUBROUTINE RETURN
A-THE (M+1)TH COLUMN CONTAINS THE APPROXIMATING VECTOR AB
6--COEFFICIENTS OF THE FIT
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20/33/39
                                                                                                                                                        KEUCLE THE LEAST SEUAKES MATRIX TO UPPER TRIANGULAR FORM
 UATE = 79206
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                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                 ENTRY ALSUITA, Y.B. KZ., NN, MM, NA.J. U.D. U. 1=1,N.A.L. | X.L. 
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                              A(1,1)=A(1,1) - A(1,1)+A(N1,1)
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                             UC 50 1=L,N
PF = PP +A(1,L)*A(1,M1)
U=PP/(-A(L,L)*A(N1,L))
UC 100 1=L,N
A(1,M1)= A(1,M1) - U*A(1,L)
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                               ALSO
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                             B(M)=A(M,M1)/A(N1,M)
1F(M-EQ-1) GC TO 130
UU 120 LL=1,MM1
L=M-LL
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PF=PP-A(L,1)+B(1)
O(L)=FP/A(N1,L)
                                                                                                                                                                                                                                          UC 10 1=L.N
55=55+A(1)-L1**2
                                                                                                                                                                                                                                                                                                             S=05ER1(52)
1F(a(L).L).L1.0
0=52 + 5*A(L).L
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IF(L, EQ.M) 6C
L]=L+1
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26/33/39
                      SS=0.bc
MFI=M+1
UC 14C 1=MFIPN
SS=SS+4(1,PM) **Z
A(1,M) =0.bd
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ALSW CALLED, TIME 0.4100

RESIDUAL SUM CF SCUARES 0.39550500060570-08

RESIDUAL SUM CF SCUARES C.1711446421510-08

RESIDUAL SUM GF SCUARES 0.1711718754270-08

COEFFICIENTS FUR 1 CELLS CUMPUTED, DELTATE 0.

.TAT= C.1200 AVERAGE TIME= 0.1200

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THIS PALEXAM ACCEPTS CARD INPUT WHICH SPECIFIES THE MINIMUN AND MAXIMUM LIMITS OF A 1621 REGICE WHICH GUERLAPS OR IS CONTAINED IN THE SEGION ESTABLISHED AND MUDELLED IN SECTION II. THE ENTIRE REGION CONSTITUTES A SINGLE CELL WHICH IS AGAIN DIVIDED BY A REGULAR GRID OF CONSTITUTES A SINGLE CELL WHICH IS A AGAIN DIVIDED BY A REGULAR GRID OF CONTAINED BY A REGULAR GRID OF CONTAINED BY SINGLE CALCULATED USING MASS. MODEL 31G (SUBRECHINE PIMASS) AND THE MODEL SIDE OF USING THE MODEL SIDE OF USING THE MODEL SIDE OF USING THE MODEL SIDE OF SECTION A RAYDUM ACCESS FILE (FILE I). AN ERROR ANALYSIS IS ANALYSIS IN AND THE RESULTS ARE CUITOT. IN ADDITION OF ACT OF SECTION AND MARYSIS IN A MAXIMUM ADSULUTE ERRURS ARE UUIPUT ALUNG MITH THE LUTAL.
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FURMAI(1h0,1CX,y(1h-),i)(cX,c(1h-)))

FURMAI(1h0,1CX,cHr)Mass (1ME/exeCullON=,ry.c)

LO FORMAI(1h0,1CX,cHr)Mass (1ME/exeCullON=,ry.c)

LO FORMAI(1X,-5r)Fin.3,c(CFI)0,4),y(1X,5rfy.4))

LO FORMAI(1X,1y+simmern)

LO FORMAI(1X,
UAIL = 79206
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CUMMUN / FLDATA/KEA(Y),IN7(5)
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AIF 1M1=K1F1S

ESC=LCO/K1F1S

ESC=LCO/K1F1S

ESC=LCO/K1F1S

ESC=LCO/K1F1S

ESC=LCO/K1F1S

ESC=LCO/K1F1S

SLEGU=DSCK1 (SOECE/K1F1M1)

SUECE=DSCK1 (SOECE/K1F1M1)
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stebe=stebu+lnkbu+erkbu
stebn=stebn+lnkbn*ekkbn
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MCDIFICATIONS 3Y JUHN L. JUNKINS, KEMI C. ENGELS, AND JUHN J. SMITH DEPAKIMENT DE ENGINEEKING SCIENCE AND MECHANICS SCHOOL UP ENGINEEKING VIKGINIA PÜLYTECHNIC INSTITUTE AND STATE UNIVERSITY ELACKSBURG, VIKGINIA 24061

APKIL 111777 OF LAST MUDIFICATION -

URSIARU FUR P. SOME PULNT IN SPACE
X,Y,Z—EAKIM—FIXED KECTANGULAK COOKDINATES
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NURDER——UKUER UF INE BASIS FUNCTIONS DESIRE

INFOIN

HMIN, ALMIN, APMINIMUM BOUNDS OF FINITE ELMENT FIELD DESIRED NMAX, ALMAX, APMAX, APMAXIMUM BOUNDS OF FINITE ELEMENT FIELD DESIRED MCELL, ALCELL, APCELL --CELL SIZE IN M, LAMBOA, PHI NHAN, NFHI --NOMBER OF LELLS IN M, LAMLA, PHI DIRECTION NOCKOCK --LKOER OF FULYNOMIALS DESIRED NCHOCK -- LKOER OF FULYNOMIALS DESIRED NCHOCK -- LKOER OF FULYNOMIALS DESIRED NCHOMBER OF COEFFICIENIS IN THE MODELING EQUATION TRUM LUMMUN FLUATA

LISK

CUICE, CN-MUDICING CUEFFICIENTS FUR THE HILAMDA, PHI CUMPUNENTS INE GRAVIITY ANCHALY FREM CUMMUN NEW TERMUN NEW TERMUN NEW HILL FROM NORMAL ID PHINTELED IN A ALAMINIA ASOVE REF. ELLIPSUID ALUNG NORMAL ID PHINTELED IN A ALAMINIA (KAUIANS) OF ALAMINIA FEUDE ILCOEDCENIAL LENGIIUUE (KAUIANS) OF PHINTELED IN GEORGE ICC LAITTUDE (KAUIANS) OF P

5

7

PRUCESS

THIS PRECKAM ACCEPTS INPUT OF ELLIPSGIDAL COORDINATES (H.LAMDA, PHI)
FROM THE MAIN PROCKAM FINIES AND EVALUATES THE THREE GRAVITY
ANDMALIES AT THIS FOINT.
FIRST THE THEOLOGY TO SECTION IS TESTED TO SEC IT II
LIES WITHIN THE FEGIUM MUDELED IN SECTION IS TO NOT! THE CRAVITY
DISTURBANCES AND A MESSAGE IS PRINIED INDICATINE FINASS WHICH USES MASS
MOLE SIGN AND A MESSAGE IS PRINIED INDICATINE FINASS WHICH USES MASS
CONTAINS THIS FOINT IS DETERMINED AND THE REGION) WHICH CONTAINS THIS FOUNT IS DETERMINED AND THE AVAILABLE IN CONCERNING SEE IF THE MODELING COFFICIENTS FOR THIS CELL ARE AVAILABLE IN CORDINATES
AND SECTION WHICH THE COLFFICIENTS ARE AVAILABLE IN CONCURNINATES
AND SECTION OF THE POLITICIANTS ARE AVAILABLE IN THIS PLAINT
FOUNT TO THE POLITICIANTS AND THE SPECIFIED CROSEN.
FOUNT TO THE POLITICIANTS AND THE SPECIFIED CROSEN.
FOUNT TO THE POLITICIANT OF THE SPECIFIED CROSEN.
FOUNT TO THE POLITICIANT OF THE BASIS FUNCTIONS OF PREDETER. 

FORTRAN IV GI	KELEASE Z.O. FINITE C. MINED URGERS. IHIS GIVES
	C EACH OF THE THREE GRAVITY ANDMALIES. WHEN THE THREE CONTRIBUTIONS C OF ALL THE CLEFFILLENTS ARE SUMMED SEPARATELY. THE RESULTS ARE THE C H.LAMDA, Ph. COMPONENTS OF THE GRAVITY ANGMALY AT THE GIVEN POINT. C THESE COMPONENTS ARE RETURNED TO THE MAIN PROGRAM FINITS.
	LUIPUIS
	C 10 CUMMUN HLP C 60,64,6NELLIPPUIDAL COMPONENTS OF THE GRAVITY DISTORBANCE
2000	IMPLICA
0.000	COMMON VALYANTALAMANTALAGO GENGRALAGO COMMON VALYANAYA GANGA GANGA COMMON VALAGO SANDANAMANA MALELA ALMINA ALMAXA ALCELLA APMAXA COMMON VALUA AMANA HILELA ALMINA ALMAXA ALCELLA APMAXA
2000	C SIMENSION IX(7), TY(1), 12(7), CC(5+), CE(6+), CN(5+)
JOCK	LIMENSION IXIO+1, IYIO4), IXIO4)
4000	UAIR JA/ 1919-1919 1919-1919-1919-1919-1919-191
0100	OAIA 18/10/10/10/10/10/10/10/10/10/10/10/10/10/
	*5969 (91969594959691969594959196959491969517 UAIA 12/11/11/11/11/11/11/11/11/11/11/11/11/1
2100	5
	C CLEMENT CLURATINATES TO REGION BUNNUARIES. AF PUINT IS COTSIDE FINITE
6013	** IFI(Hallamin).Un. (Has).Uk. (ALAM.LI.ALMin).Uk. (ALAM.C).ALMAX *** UK. (APHI-LI.APMIN).Uk. (APHI.C).APMAX)
	C CLMPUTE VARIOUS INDEXES TO HELP FIND THE KIGHT CELL IN THE KEGICN
50015	-17K
84000 0000 0000	IT (TOELONGAX) A=1-1 IF (ALAMORE OF ALMAX) C=C-1 IF (APPINET OF AND N=K-1 X1-1
00.23 00.23 00.23	
	C CALCULATE THE RECORD PUINTER FOR THE SET OF COLFFICIENTS AT THESE C CUCRESINATES

Comment

in.

Common of the last

20/24/03 SINC=USIN(ALAM)

CUSP=UCOS(AFM1)

SINC=USIN(AFM1)

SINC=USIN(AFM1)

SINC=USIN(AFM1)

GE=COS(AFM1)

GE=COS(AFM1)

GE=COS(AFM1)

GE=COS(AFM1)

FUSIN(AFM1)

FUSIN(AFM1)

FURMAT(49M FUIN) NUT IN FINITE ELEMENT FIELD, PIMASS CALLED)

KEI IN=UMN

KEI UMN UATE = 79206 FINITE FURTRAN IV G1 RELEASE 2.0 

FUNTRAN IV 61 MELEASE 2.0

20 TA(1)=2.\*XBAK\*TA(1-1)-TA(1-2)
ENURN

DATE = 7920c

20/34/03

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PA
                                   20/34/03
                                   UATE = 79206
                                                IF (PMVALS(1)) 10,20,10
CONTINUE
DX=PUSITS(1,1)-X
UY=PUSITS(1,4)-Y
UZ=PGS113(1,4)-Y
UIST=USH1(1,5)-Z
UIST=USH1(D1STSH)
TEMP=PMVALS(1)/UIST/D1STSH
                                 PIMASS
                                                                                                               DELCX=DELGX+(DX+1EMP)
LetcY=DelcY+(DY+1EMP)
DelcZ=DElcZ+(GZ+1EMP)
in cuminde
                                                                                                                                                          KETURN
ENU
                                   RELEASE 2.0
                                                         2
                                FORTRAN IV G.
                                                 $00000000 0000 3
                                                                                                                                                        C021
0022
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E S S S S S S S S S S S S S S S S S S S	LAMBUA	int G	ENNED (MCALS)	LAKEL	(MCALS)	FMGU (MGAL >)	PMGE	IMCALS)	F160	(MOALS)	ולוטעראו
13	10	1.	0.37.0	0.0000	0.0101	-0.4254		1	-0.6180	1	(1,0063
3	212	3,4	4771-0	1.3/54	2+77-0	-4.9000	ℳ.		-3-0210	4.5	7.3070
33	191	:;	0007.0	-1.1554	0-1823	-2.3644	, ,		-4.3554	::	-0.1633
3			10300	0187-0	0.0330	14.7033	٠.	•••	-4.3603	700	2000
3			****	20110	-1.1649	-6.0700	, ,		10000	3.6	1.1401
3		.,	1.6151	-0.3063	1701-1-	14.3715	A 4	-	-5.79.00		-1.5760
0000	75.6333	-24.6333	1114.1-	1477	7555	-12.1221	,	**	-10-5116	3.5	4.5 2 30
3			-0.400	0.0254	-1-696	-11.8045	4	~	-10.83/8	3.4	1.6166
33			0.407.0	-0.5360	1.0738	14.36.4-	v -	: :	-10-1334	7.7	7.5/63
3		7.0	-1.1101	-0.6454	0.0646	-14.4645	-		-13.3644	7.7	2.1011
35	ns		0.0048	20000	-0-2001	-13.7666	-	:-	-14-3634	200	-1-1-1-1
3	1.		-0.0144	-0.6367	-0-4066	-0.8464	v.	1	-0.6140	4	2.5140
33	U .		2000	10.000	000000	11.0000	<i>J</i>		10.413	77	1.1.70
3	131	:	-1.6400	10.545.0	C . 30 8 3	1501-7-	0		-1.5565	:-	-1.3273
3	414		2.2843	7407-0-	** 11.0-	74.0960	4"		-4.4612		5521 - 7
33		3.0	757000	10.46.48	0.0400	1401-5-		-	-3.86/1	•	0.0148
3	so.	-	-0-4154	4T40-0-	0.240	-3.4055	-	7	1056-7-	7	-1.144
33			10000	2270	-0-1212	-3.3678	2 S	1 0	-3.1112	, 4	301921
3	1 11		0.5271	-4.3664	0-7126	-4.57+Y			-5.5660	2	
3	S	-	-0.2161	01750	1.0693	-4. (433	<b>T</b>	7	-4.5373	-	-1-4-1-
			10000	0.4380	10000	19.4081	T J		00000	: :	200000
3			1.2043	-1.1760	6.4263	4874-4-	•		-1.1934		01410
200		-,	0.25.71	20000	0.1085	-6-1014 -1-1-1-1-1-1-1-1-1-1-1-1-1-1-1-1-1-1-	V		-6.3910	2-	-1-4050
3	12		0.4411	6667-0-	-0.0402	6067-1-	, _	'n	****		C.5 (81
37.00	J.		0.5179	-0.08AC	-0-01de	-1.4335	2	7	-1-1514	3	0.6339
22	7.7		0.3%	10.00	0.414	-1.5656	NV	42	1.9/45		700000
3	n.	3.	0.5431	-6.2241	50.00	-1-6284		v.	-4.3640	1	0.1640
00.00	תח		10.5651	77770	0-00-0	-2.0/30			-1-1-1485	, ,	-6.3330
30-04		7.	0.1512	0.0141	0.1043	-4.1845			-4.3351	4.	0.0150
		7.0	0100	10.000	0-62-0	7995	nv		1696		10.10
90.00	n	7.	1764-0-	-0-140	0-1037	15.6450			-4.1944		BOS 5.0-
	4 21		7714-0-	207.0	6.2020	-20031	7		-3.4 (45	77	0.17.00
BO.CC			0.6316	1.01.0	C. 1019	-3-1785		3	-3.4100	:	
20.00			-0.054	224-0-1	2040-0-	13.6181	-		-4.7593	2.4	10.4.01
2	12	4.6		-0.4470	1617-0-	-1.6345		: -:	-0.9104	3.5	6.3003
30-37	•	5.3	271100	-0.3039	6.0000	-1.0910	•	-	-U. VIVB	3	
	.,,,,		0.4440	2000	87.77	-1.6133	$\mathbf{u}$		11.144	3	414000
וניינו	-11	7.0	-1.1154	-1-1 125	D. 22.0-	-1.3025	n.	7	-1.4873		(144.0)
	, .,		10.17.7	11111	1757	1.00001			11.1544		
30-01	•	2.	2448.0	-6.6873	-6.1920	9255-1-	0		41.69.1-	2.	2000
	.,,,	2	10.530	2012	201000	11.06.01	ש ס	-	1100011	97	16717
00-01	311	7.	-0.3533	-0-105	-0.0665	3770-1-			-1.4695		-6.3465
		7.5	177777	100000	-0-1502	1764-1-	05		-1-0134	4 5	550000
30-01			-0-1145	(12.153	0.0344	-1.96.15	·	7	-1.1350	0	10.1140
33.01	*1		-(-115/	140000	-0.1e55	-1.66%	D	•	-1.1135	٥.	-1-6425
724			70								
ABS. VALUES	IN TALNE	KKCHS	1.3111	1.3626	1.5247						
	-	VI ALLUINA	, , ,								
•	IMASS TIME!	/ Extending	10504701 =								

ORTHONORMAL POLYNOMIAL FINITE ELEMENT SOFTWARE

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WRITE(6,4) NORDER,NC
WRITE(6,4) MGBSU,MGBSE,MGBSN,MOBS
WRITE(6,5) HCELL,ALCELL,APCELL,HMIN,ALMIN,APMIN,HMAX,ALMAX,APMAX
22/16/52
                                                                                                                                                                                                                                                                                                                                                                                                        GET THE COEPFICIENTS OF THE CRIMOGONIAL POLYNOMIALS FOR A TYPICAL
CELL
                                                                                                                                                             COMPUTE THE NUMBER OF FINAL CELLS IN UP, EASTERN, AND NORTHERN DIRECTION AND THE TOTAL NUMBER OF FINAL CELLS IN THE FINITE ELEMENT FIELD
= 79206
                                                                                                                                                                                                                                                                                              SIZE WITHIN EACH CELL.
DATE
                                                                CCNVERT ANGLES FROM DEGREES TO RADIANS
                                                                                                                                                                                                                                                                                                                                                                                                                                     CALL CRIEC(P1, NORDER, MUBSU, DIH)
X1=-1,00
DO 12C IH=1, MUBSU
CALL MULT(X1, NORDER, P1, TX(1, IH))
X1=X1NUE
CGONTINUE
CALL CRIEC(P2, NORDER, MGESE, DIL)
X2=-1,00
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                         16 146 Ir=1,MOuSN
ALL MUL! (X3;NUKDER,P3,12(1,1P))
3=X3+D1P
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                          130 130 1L=1,MUBSE

ALL MULI(XZ,NCKDER,PZ,TY(1,1L)

(2=x2+01L

CNTINUE

ALL UAIND(P3,NURDER,MUESN,DIP)

3=-1.00
                                                                                                                                                                                                        RNH=(HMAX-hMIN)/HCELL+.59900
RNL=(ALMAX-ALMIN)/ALCELL+.79900
RNP=(APMAX-APMIN)/ALCELL+.79900
NH=RND
NLAM=RNL
NPHI=KNP
                                                                                   ALCELL=ALCELL#DEGRAD
APCELL=APCELL#DEGRAD
ALMIN=ALMIN*DEGRAD
APMIN=APMIN*DEGRAD
ALMAX=ALMAX*DEGRAD
APMAX=APMAX*DEGRAD
MAIN
                                                                                                                                                                                                                                                                           NCELLU-NI+NLAM+NPHI
                                                                                                                                                                                                                                                                                              COMPUTE THE INCKEMENT
                                                                                                                                                                                                                                                                                                                   DH=HCELL/RMCM1
DLAM=ALCELL/RMEM1
UPH1=AFCELL/RMNM1
                                                                                                                                                                                                                                                                                                                                                              DIH=Z.DOZKMUM]
DIL=Z.UOZKMEM]
DIP=Z.DOZKMRM]
2.0
RELEASE
                                                                                                                                                                                                                                                                                                                    110
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                             120
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                             130
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FCRTRAN IV G1
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FORTRAN IV G	G1 RELEASE 2.0 MAIN DATE = 79206 22/16/52
	C NUTE: IF A SYMMETRIC GRID PATTERN IS ALWAYS CHOSEN (SUCH AS 4***4) C THE THREE SETS OF EVALUATIONS ABOVE ARE REDUNDANT AND TWO COULD BE
1000	C WRITE FIELD DATA ONTO FIRST RECOND OF SEQUENTIAL FILE
0076	WRITE(3) hMin, hMAX, hCELL, ALMIN, ALMAX, ALCELL, APMIN, APMAX, APCELL, PI, +PZ, P3, NH, NLAM, NPHI, NUKDEK, NC, MUBS, NCELLS
	L FILL THE A MATRIX
72.000 72.000 73.0000 73.000 73.000 73.000 73.000 73.000 73.000 73.000 73.000 73.0000 73.000 73.000 73.000 73.000 73.000 73.000 73.000 73.000 73.0000 73.000 73.000 73.000 73.000 73.000 73.000 73.000 73.000 73.0000 73.000 73.000 73.000 73.000 73.000 73.000 73.000 73.000 73.0000 73.000 73.000 73.000 73.000 73.000 73.000 73.000 73.000 73.0000 73.000 73.000 73.000 73.000 73.000 73.000 73.000 73.000 73.0000 73.000 73.000 73.000 73.000 73.000 73.000 73.000 73.000 73.0000 73.000 73.000 73.000 73.000 73.000 73.000 73.000 73.000 73.0000 73.000 73.000 73.000 73.000 73.000 73.000 73.000 73.000 73.0000 73.000 73.000 73.000 73.000 73.000 73.000 73.000 73.000 73.0000 73.000 73.000 73.000 73.000 73.000 73.000 73.000 73.000 73.0000 73.000 73.00000 73.0000 73.0000 73.0000 73.0000 73.0000 73.0000 73.0000 73.00000 73.0000 73.0000 73.0000 73.0000 73.0000 73.0000 73.0000 73.00000 73.00000 73.00000 73.00000 73.00000 73.00000 73.00000 73.000000 73.0000000000	DC 150 11=1,NC NX=1X(11) NY=1Y(11) NZ=1L(11)
OCC OC	DC 15C 1H=1,MGESU DC 15C 1L=1,MCESU DC 15C 1L=1,MCESU
200	
2000	150
	C INCREMENT THE H COORDINATE (HMCELL) I CELL SIZE
	C HMCELL=HMCELL+HCELL
	C RESET THE LAMBA COURDINATE TO 115 CELL MINIMUM VALUE
0092	ALMCEL=ALMIN-ALCELL DO 30C ILC=1, NLAM
	C INCREMENT THE LAMBA COORDINATE (ALMCEL) I CELL SIZE
*****	ALMCEL+ALCELL
	C RESET THE PHI CUGRUINATE TO ITS CELL MINIMUM VALUE
9600	APMCEL=APMIN-APCELL DO 3CC IPC=1,NPHI
	C INCREMENT THE FMI COURDINATE (APMCEL) I CELL SIZE
0047 00348	APMCEL=APMCEL+APCELL MJ=C
4400	11

RELEASE 2.3 MAIN DATE = 79206 22/16/52 DO 250 IH=1,MOBSU	N.	ALAM-ALMCEL-ULAM DC 250 IL=1,MGBSE	C INCREMENT THE LAMBA COORDINATE (ALMCEL) I GRID INCREMENT SIZE	ALAM=DLAM+DLAM COSL=DCOS(ALAM) SINL=CSIN(ALAM) APHI=PMCEL-UPHI DC 25C IP=1,MOBSN	C INCREMENT THE PHI COORDINATE (APMCEL) I GRID INCREMENT SIZE	APHIMPHI-UPHI CCST-CCUS(ATHI) SIRT-SSIN(ATHI)	C TRANSFURM ALL ELLIPSUIDAL COURDINATES INTO X,Y,Z COORDINATES	C RN=DSCKT (A2*CCSP+B2*SINP*SINP)  LN=B2/KN  KN=A2/KN  Y=(KN+H)*CLSP  X=Y*CUSL  Y=Y*SINL  Z=(2N+H)*SINP	C DETAIN THE GRAVITY CESERVATION AT THIS GRID PUINT	CALL PIMASS	C DIRECTION CUSINE MATRIX FUR 3-2 RUTATION (LAMBDA,-PHI) FROM X,Y,Z, 10	C C11=COSP*COSL C C12=CUSF*SINL C C13=SINL C C2 = -SINL C C2 = -SINL	C. S.	GU=CLSF*CDSL*GX + LOSP*SINL*GY + SINP*6Z GE=-SINL*GX + CUSF*GZ GN=-LNP*CUSL*GX + CUSP*GZ GN=-LNP*CUSL*GX + CUSP*GZ	C STORE THE UBSERVATION DISTURBANCES
O100		0101		00000		6109 61110 61112		00113 00115 0117 0119		0120				0122	SCHEENING DESCRIPTION

PA

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22/16/52
                                                                    WRITE THE THREE COSERVATION ARRAYS UNTO THE SEQUENTIAL FILE
  DATE = 79206
                                                                                   (DELGU(I), 1=1, MOBS)
(DELGE(I), 1=1, MOBS)
(UELGN(I), 1=1, MOBS)
  MAIN
                      200 DELGU (MJ) = GU
DELGE (MJ) = GE
DELGA (MJ) = GN
250 CON 11 NOE
                                                                                  WARITE (3)
WARITE (3)
CONTENS
END INCE
RELEASE 2.0
                                                                                                         300
                                                      JUUU
FORTRAN IV G1
                      0125
                                                                                  00128
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FORTRAN IV G1	RELEASE 2.0 GRTHO DATE = 79206 22/16/52 SUBROUTINE URTHUIP, NURDER, MUB, DELTAL)
	DEPARIMENT OF ENGINEERING SCIENCE AND MECHANICS SCHOOL OF ENGINEERING VIRGINIA POLYTECHNIC INSTITUTE AND STATE UNIVERSITY BLACKSBURG, VIRGINIA 24001
	C DATE OF LAST MUDIFICATION APRIL 1,1979
	L INPUT
	C FROM SUBROUTINE CALL C NCFDER PRDER OF BASIS FUNCTIONS DESIRED C MCGNUMBER UF INTERVALS IN THE FINITE ELEMENT CELL C DELTALSIZE OF FACH INTERVAL
	C PRUCESS
	C THIS SUBKCUTINE ACCEPTS THE BASIS FUNCTION ORDER, THE INTERVAL SIZE OF THE FINITE ELEMENT CELL TO BE MODELED.  C AFTER THE MATRICES ARE INITIALIZED, THE NORMALIZED INNER PRODUCTS AND CFINAL CCEPFICIENTS ARE CALCULATED. SINCE THE FIRST EQUATION, URDER 1 C. AND CROEK 2, FZ=X, ARE ANCWN, CNLY THE INNER PRODUCTS ARE CALCULATED. SUNCE THE FIRST EQUATION.
	Coutput
	C TO SUBRCUTINE RETURN C PTHE INNER PRODUCTS USED TO CALCULATE THE BASIS FUNCTIONS VALUES C RECURSIVELY
20003	1
8	C INITIALIZE THE MAIRICES THAT WILL HOLD THE FINAL CUEFFICIENTS OF
0000 0000 0000	
55	F(J, I)=0.06 20 CCNTINUE C USING THE 'EVEN' PROPERTY OF THE URTHUGUNAL EQUATIONS, CALCULATE THE C SUMS: BLP=1+X2**I+X.2**I+XN**I, WHERE I=2,3,4,*CUSS, AND C SUMS: X=THE NORMALIZED COORDINATE OF ACH GRID POINT

```
CALCULATE THE CCEFFICIENTS OF THE URTHOGONAL EQUATIONS OF DRUEN I BY USING THE RECURSIVE PROCESS: F(I+1)=F(I)*X-P(I)/P(I-1)*F(I-1)
WHERE P(I) IS THE INNEW PRODUCT OF F(I)*F(I)

DU +00 I=1.N
 22/16/52
                                                                                                                                                                                                                                                                                    COEFFICIENTS OF F(1+1).
                                                                                                                                                                                                                                                                                                                                                300 CONTINUE
40C CONTINUE
NORMALIZE THE COEFFICIENTS OF THE ORTHOGONAL POLYNOMIALS F.
DC 6CO J=1,N
P(J,1)=DSCRI(P(J,1))
P(J,1)=DSCRI(P(J,1))
DC 6CO [=1,N,201,1)
 DATE = 79206
                                                                                                                                                                                                                                                                                   CALCULATE THE
                                                                                                                                                                                                                              P(1,1)=P(1,1)+F(K,1)+F(L,I)+BLP(LP)
CCN1 inue
                                            XL=-1.500-06-11AL

XL=XL+DE-14MUS

XL=XL+DE-1AL

IF (DABS(XL)-Le.1.0-15) GO TO 100

3LP(I)=BLP(I)+XL**J
                                                                                                                                                                                                                                          JF((1,-LE-1)-CK-(1,-EE-N) 60 10
                                                                                                                                                                         (1,1)=1.00
ULATE THE INNER PRODUCT P(1)
U 200 K=1,1
C 20C L=1,1
P=K+L-1
                                                                                                                                                                                                                                                                                                                                         (L,K)=F(M,1)-P(J,2)*F(L,J)
ONTINUE
                                                                                                                                                                                                                                                                         K=1+1

NG THE INNEK PRODUCTS C

P(J,Z)=P(J,1)/P(J,1)

F(1,K)=-P(J,Z)*F(1,J)

DC JCC L=2,J
DRIND
                   BLP(1)=MOB
DG 1G0 1=2,MOES
J=1-1
BLP(1)=0.00
XL=-1.00-06L1AL
XL=-XL+0FL1AL
 RELEASE
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FORTRAN IV G1
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A
 22/16/52
DATE = 79206
              G CONTINUE

UX=PUSITS(1):1) -X

UX=PUSITS(1):2) -Y

UZ=FUSITS(1):2) -Y

UISTSC=UX*UX*UY*UZ*UZ

UISTSC=UX*UX*UX*UZ*UZ

UISTSC=UX*UX*UX*UZ*UZ

UISTSC=UX*UX*UX*UZ*UZ
PTMASS
                                                                                      DELGX=DELGX+(DX*TEMP)
DELGY=U=LGY+(LY*TEMP)
DELGZ=DELGZ+(02*TEMP)
                                                                                                                      20 CUNITNUE
                                                                                                                                      RETURN
RELEASE 2.0
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                                                                               J
FURTRAN IV G1
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                                                                                                                     0050
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DA

-SHOO DEPARTMENT OF ENGINEERING SCIENCE AND MECHANICS SCHOOL OF ENGINEERING VIRGINIA POLYTECHNIC INSTITUTE AND STATE UNIVERSITY BLACKSBURG, VIRGINIA 24061 JOHN L. JUNKINS, REMI C. ENGELS, AND

LAST MODIFICATION -- APRIL 1,1979 9 DATE

INPUTS

FROM DISK
FILE 3
HMIN, ALMIN, APMIN--MINIMUM BOUNDS OF FINITE ELMENT FIELD DESIRED
HMAX, ALMAX, AFMAX--MAXIMUM BOUNDS OF FINITE ELEMENT FIELD DESIRED
HMAX, ALMAX, AFMAX--MAXIMUM BOUNDS OF FINITE ELEMENT FIELD DESIRED
FILE, ALCELL, ACCELL --CELL SIZE IN H, LAMBOA, PHI
A - ACH DIRECTION
IN - ACH DIRECTION
IN - ACH DIRECTION
NORDER--CEUER OF CELLS IN H, LAMBA, PHI DIRECTION
NORDER--CEUER OF POLYNOMIALS DESIRED
NOC--NOMBER OF COEFFICIENTS IN THE MODELING ELUATION
NORDER--IHE ICIAL NOMBER OF CELLS IN THE MODELING AREA
A-IHE LEASI SCUARE MAIRIX
DELGU, DELGE, DELGN--THE GRAVITY DISTURGANCE OBSERVATIONS IN EACH
COURDINATE DIRECTION

PROCESS

THIS PROGRAM ACCEPTS AS INPUT DATA WHICH DESCRIBES THE REGION TO BE MUCELED AND THE CONFILLENTS OF THE ORIHHNDRMAL BASIS FUNCTIONS, FROM DISK FILE 3. THIS INFORMATION IS IN THE LEAST SQUARES MATRIX A. IS UNE THE SANDOM ACCESS FILE (FILE 1). THE LEAST SQUARES MATRIX A. IS FROM FILE 3. NEXT BELL STONES OF FILE 10. THE LEAST SQUARES MATRIX IS FROM FILE 3. NOW THE LEAST SQUARES MATRIX IS FROM THE LEAST SQUARES MATRIX IS FROM THE LEAST SQUARES MATRIX BY THE DISTURBANCE VECTOR TO PRODUCT THE MODELLEAST SQUARES MATRIX BY THE DISTURBANCE VECTOR TO PRODUCT THE MODELLEAST SQUARES MATRIX BY THE DISTURBANCE VECTOR TO PRODUCT THE MODELLEAST SQUARES MATRIX BY THE DISTURBANCE VECTOR TO PRODUCT THE MODELLEAST SQUARES MATRIX BY THE DISTURBANCE OF FILE IS INVOLVED THE COEFFICIENTS ARE WRITTEN IN THE NORMAL UNDER CELL IS INVOLVED THE COEFFICIENTS ARE WRITTEN IN THE NORMAL UNDER CELL IS INVOLVED THE COEFFICIENTS ARE WRITTEN IN THE NORMAL UNDER THEN VALLES MOST RAFIDLY, FOLLOWED BY LAMDA AND FINALLY H. 

TO DISK

BOUNDS OF FINITE ELMENT FIELD DESIKED BOUNDS OF FINITE ELEMENT FIELD DESIKE SIZE IN H.LAMSDA,PHI HMIN, ALMIN, APMIN——MINIMUM HMAX, ALMAX, APMAX——MAXIMUM HCELL, ALCELL, APCELL——CELL 22/17/21

1

- Comment

C.0700 AVERAGE TIME= 1 CELLS COMPUTED, DELTAT= 0.1900 ALSO CALLED, TIME= COEFFICIENTS FOR

0.0700

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THIS PROCKAM ACCEPTS CARD INPUT WHICH SPECIFIES THE MINIMUN AND MAXIMUM LIMITS OF A TEST REGION WHICH OVERLAPS OR IS CONTAINED IN THE REGION ESTABLISHED AND MUDELLED IN SECTION 11. THE ENTIRE REGION CONSTITUTES A SINGLE CELL WHICH IS AGAIN DIVIDED BY A REGULAR GRID DISTURBANCES ARE CALCULAR GRID BY A REGULAR GRID DISTURBANCES ARE CALCULATED BY USING THE MODELLED GRAVITY DISTURBANCES ARE CALCULATED BY USING THE MAKE DETERMINED IN SECTION IN AND STORED ON EACH TEST FOINT AND THE RESULTS ARE USING ADDITIONS.

THEN TO THE PUBLIC ERRORS ARE USING WITH THE TOTAL OFFICE AND MAXIMUM ADSCLUTE ERRORS ARE USING WITH THE TOTAL OFFICE OF THE TWO METHODS.
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                          FORMAI(F10.0,2F10.2)
FORMAI(F10.0,2F10.2)
FURMAI(1X,12110)
FURMAI(1X,F12.0,2F10.2)
FORMAI(1X,F12.0,2F10.2)
FORMAI(1R0,4A,1HH,6X,0HLAMCJA,5X,3HPH1,7X,5HEKKGU,5X,5HEKKGE,5X
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                               7 FLRMAI(4X,4H(KM),1X,2(5X,5H(DEG)),2X,9(3X,7H(MGALS)))
8 FLRMAI(2X,9(1H-),11(2X,8(1H-)))
9 FLRMAI(1HC,1CX,2ZHFINIE TIME/EXECUTION=,Fy.c)
10 FURMAI(1HC,10X,2ZHFINIE TIME/EXECUTION=,Fy.c)
11 FORMAI(11X,-3FFINO,4,2TCFINO,+),9(1X,5FFY,+))
12 FORMAI(1X,-3FFINO,4,2TCFINO,+),9(1X,5FY,+))
13 FORMAI(1X,-2H-SHMCAND,-2X,3(1X,5FY,+))
14 FORMAI(1X,-ZCHAS), VALUES OF MAX. ENRORS,4X,3(1X,5PFY,+))
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                            IMPLICIT KEAL*E (A-H,O-Z)
COMMUN /HLP/H,ALAM,APHI,GU,GE,GN
COMMUN /XYZ/X,Y,Z,GX,GY,GZ
CGMMUN/IMAKK/IFLAG,1S1IIN
COMMUN/FLDATA, REA(9),REB(14),REC(14),REU(14),INI(2)
COMMUN /MASPUS/PMVALS(100),PUSITS(1080,3)
                                                                            HMAX, ALMAX, APMAX -- MAXIMUM BOUNDS OF TEST FIELD DESIRED
   DATE = 79206
                                                                                                                                                  THE CARD INPUT FOR THIS SAMPLE RUN IS AS FOLLOWS:
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                             DEFINE FILE 11251, COIG, U, IPUINT)
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                      DATA A/6376160.00/
DATA B/0356774.50400/
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                              ERRUR ANALYSIS REPURT
                                                                                                                                                                                                                                                                                                              PAOCESS.
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FURTRAN	1 IV 61	RELEASE	SE 2.0 MAIN DATE = 79206 22/17/43
0024		٠	15 FORMATICACHO NEGATIVE UR ZERU NU. UF STEPS)
		.د.ر ر	READ IN THE FIELD DATA FROM THE FIRST RECORD OF THE RANDOM ACCESS FILE (FILE 1).
9700		. د	READ(1-1) REA, MED, MEC, MED, IN]
		ى د د	READ FILE & WHICH CONTAINS THE MASS VALUES AND THEIR X,Y,Z COOKGINAT SET UP CONVERSION FACTORS AND INITIALIZATION
9787017 9897070 9000000		, د	REAU(2) PMVALS, PUS, IS PI=JAKCUS(-1.DU) DEGRAL=PI/180.DU A2=A4A B2=E48 IFLAG=C ISTIN=U
		ررر	KEAD IN THE TEST CELL GRID PAITERN AND THE FINITE ELEMENT FIELD SCUNDS
0000000 0000000 0000000000000000000000		, ,	CU READ(2,1,enD=4C) INTERMINISTERLISTERN READ(5,2,enD=4C) hMIN,ALMIN,APMIN,HMAX,ALMAX,APMAX WAITE(6,3) INTERMINISTERLISTERN ARITE(6,4) INTERMINISTERN INMITE(6,5) HMIN,ALMIN,APMIN,HMAX,ALMAX,AFMAX INMITE(6,5) HMIN,ALMIN,APMIN,HMAX,ALMAX,AFMAX INMITEINTERLIT IPMI=1NTERLI
		ارور	VERIFY THAT THE TEST CELL GRID PAITERN IS VALID AND CALCULATE THE GRID INCREMENT SIZE
1111		J	21 WRITE(6,15) GCID 40 22 DH=0.b0
***			25 8HI = IHMI 08 = IHMI = IHMI 24 IF (ILMI) 21,25,26 25 DLAM=C.DC
00000000 00000000 00000000000000000000			20 RLM1=1LM1 27 IF (IPM1) 27 IF (IPM1) 21,20,25 28 DPH1=0,00 29 RPM1=1PM1 29 RPM1=1PM1
		ارد	SET UP THE ERPLIN ANALYSTS REPCAT READINGS
9000		<b>)</b>	30 WKITE(6,c) WYTE

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A.
22/17/43
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                 EVALUATE THE GRAVITY DISTURBANCES USING THE MUDELING EQUATION WHOSE CUEFFICIENTS WHERE DETERMINED IN SECTION II
                                                                                                                                                                                                 S GRAVITY DISTURBANCES AT EACH TEST GRID POINT
                                                                                                                                                                                                                                                                                                                        INCREMENT THE LAMBA CUBMIDINATE 1 GMID INCREMENT SIZE
                                                                                                                                                                                                                                                                                                                                                                                                                         INCREMENT THE PHI CLORDINALE I GRID INCREMENT SIZE
79200
                                                                                                                                                                                                                                                                                                                                                                                                                                                                               TRANSFURM THE ELLIPSOIDAL COUNDINATES INTO X,Y,Z
                                           INITITALIZE ALL TIME AND ERRUR ANALYSIS VARIABLES
                                                                                                                                                                                                                                                       INCREMENT THE IN CUCRDINATE I GRID INCREMENT SIZE
 11
DATE
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                          + BZ*SINF*SINF
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                    CCSP=DCUS(AFHI)
SINP=USIN(AFHI)
RN=USGA(AZ*CUSP*CUSP
ZN=BZ/RN
RN=AZ/RN
MAIN
                                                                                                                                                                                                                                                                                                                                               AL=AL+ULAM
ALAM=AL+UEGRAD
CGSL=DCGS(ALAM)
SINL=DSIN(ALAM)
AP=APMIN-DPHI
DC 35 K=1+1STEPP
                                                                                                                                                                                                                                                                             HEN+UH
AL=ALMIN-DLAM
JC 35 J=1,1STEPL
                                                                                                                                                                                                                        HENMIN-DH
UC 35 1=1,15TEPH
                                                                                                                                                                                                                                                                                                                                                                                                                                                            APHI = AP * DEGKAD
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                          Y=(RN+H)*COSP
X=Y*CESL
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                           7-12841 TANA 1-12
                                                                                      DETERMINE THE
                     WR1TE(6,3)
                                                                 F11Mt=0.00
                                                                                                                                                                                                                                                                                                                                                                                                                                                AP=AP+UPH1
RELEASE 2.0
                                  در
                                                                                                                                                                                                                                                CCC
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                                                                                                                                                                                                                                                                                                                                                                                                                                                                      CCC
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FORTRAN IV
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C083 0084

FCRTRAN IV G1	RELE	LEASE 2.0 MAIN		DATE = 79206	22/11
	JUJ	START THE TIMER TO DETERMINE THE EXECUTION TIME OF THE EQUATION	TERMINE THE EXECUTION	N TIME OF THE	E GUAT 1 ON
5600	، د	CALL LIMECN			
	ر د	STOP THE 11MER AND CONVERT THE 11ME TO SECONDS	WERT THE TIME TO SE	CONDS	
2600	. د	CALL TIMECK(ICP) FIIME=FIIME + ICP/IUC.UC	luc.bc		
20048 0100	, ر	GUT=60 GET=6E GNF=6N			
	ان	START TIMER TO DETERMINE THE EXECUTION TIME	INE THE EXECUTION TIN	Æ	
1010	، ر	CALL TIMEUN			
	، ر د	EVALUATE THE GRAVITY DISTURBANCES USING MASS MODEL	JETURBANCES USING MI	ASS MODEL 310	
01022	., ,	CALL FIMESS CALL TIMECKIICP) PIIME=PIIME + ICP/100.00	07:001		
	ررر	THANSPORM THE GRAVITY DISTURBANCES TO ELLIPSCIDAL CCORDINATES	DISIURBANCES TO ELL	IPSCIDAL CCCKD	INATES
0105 0106 0101	) ر	6U=COSP*COSL*6X + COSL* 6E=-SINL*6X + COSL* 6N=-SINF*COSL*6X -	GU=COSP*COSL*GX + COSP*SINL*GY + SINP*GZ GE=-SINL*GX + LOSL*GY GN=-SINP*COSL*GX - SINP*SINL*GY + COSP*GZ	752	
	ان	DETERMINE THE ERRUR VALUE BETWEEN THE TWO METHODS	ALUE BETWEEN THE TWO	METHODS	
0108 0109 0110		ERRGE = GE - CEF ERRCE = GE - CEF ERRCN= GN - CNF			
	. د	DUTPUT THE ERRCR ANALYSIS FOR THIS GRID POINT	1515 FOR THIS GRID PO	LNTC	
0111	. د	WKITE (6, 11) h, AL, A	WKITE (6, 11) H, AL, AP, ERRGU, ERRGE, ERRGN, GU, GE, GN, GUF, GEF, GNF	SU, GE, GN, GUF, G	EF, GNF
	ىد د.د	VETERMINE INE MAXIMUM AUSGLUTE ERROR VALUE FOR EACH COORDINATE DISTUREANCE	AESCLUTE ERRUR VALUE	E FUR EACH LID	RUINATE

22/11/43

FREDR FOR EACH CUCKDINATE DISTURBANCE EGUMAX=AERRGU EGEMAX=AERRGE EGNMAX=AERRGN AERRGE=DABS(ERRGU) AERRGN=DABS(ERRGN) AERRGN=DABS(ERRGN) IF(AERRGU.GI.EGUMAX) IF(AERRGE.GI.EGEMAX) IF(AERRGN.GI.EGEMAX) CALCULATE INE SUM UF THE EGU=EGU+ERRGU ددن

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22/17/43
                                                                                                                                                                                                                                                                           CUIPUT THE MEANS, STANDARD DEVIATION AND MAXIMUM ABSOLUTE ERROR
                                                                                                                                                                                                                                                                                                                                                               DUTPUT THE AVERAGE EXECUTION TIME PER PUINT FUR EACH METHUD
                                                    CALCULATE THE SUM LF THE SCUAKES OF THE EKKORS FOR EACH COGRDINATE DISTURBANCE
                                                                                                                                          CALCULATE THE MEAN AND STANDARD DEVIATION OF THE ERRURS
= 79206
DATE
                                                                                                                                                                                                                                                                                              WRITE(6,12) EGU, EGE, EGN
WRITE(6,13) SDEGU, SDEGE, SDEGN
WRITE(6,14) EGUMAX, EGEMAX, EGNMAX
FIIME=FIIME/KIPTS
                                                                                                                                                              IP1S=151EPH*151EPL*151EPP
RIPTS=1P15
RIPTM1=RIPTS-1.DO
EGU=EGU/RIPTS
EGU=EGE/RIPTS
EGU=25ENT (50EGU/FIPTM1)
SUEGU=05ENT (50EGU/FIPTM1)
SUEGE=05ENT (50ECE/NIPTM1)
SUEGE=05ENT (50ECE/NIPTM1)
                                                                                    S DEGU=S DEGU+EKKGU*EKKGU
S DEGE=S DEGE+EKKGE*EKKGE
S DEGN=S DEGN+EKKGN*EKKGN
CCNTINUE
 MAIN
                                                                                                                                                                                                                                                                                                                                                                                    WEITE(6,9) FIIME WAITE(6,10) FIIME COTO 20 COTINUE SIOP END
                     EGE=EGE+ERRGE
EGN=EGN+ERRGN
 RELEASE 2.0
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FORTRAN IV G1
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PA

+).UR.(APHI.LT.APMIN).UK.(APHI.GT.APMAX)) GJ TU +CO C COMPUTE VARIOUS INDEXES TO HELP FIND THE RIGHT CELL IN THE REGION UG15 0015 0015 0015 X=1DINI((ALAM-ALMIN)/ALCELL+1.00) K=1DINI((APHI-APMIN)/APCELL+1.00) IF(ALAM-EQ.ALMAX) I=1-1 IF(ALAM-EQ.ALMAX) J=1-1 IF(APHI.EU.APMAX) K=K-1 RJ=1 RJ=1 RJ=1
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DELLEMENT THE CONTRIBUTION THAT EACH COEFFICIENT HAS UPON THE FINAL BASIS FUNCTIONS.
        22/17/43
                                                                                                                                                                                                                                                                     HEAD I WEN SET OF CUEFFICIENTS UNLY IF PROPER SET IS NOT IN CURE
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                   EVALUALL THE BASIS FUNCTIONS AT THE TEST PUINT FUR EACH ORDER
= 75206
                                                                                                                                                                                                                                                                                                                                                                         HILAGETPOINT (CULIC), CE (IC), CN(IC), IC=1,NC)
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                    EVALUATE THE 3 CUCRDINATE GRAVITY DISTURBANCES
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                     THAN'T CHE COURDINATE KANGE (C.1) TO (-1.1)
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                      FILL THE MINIMUM CELL BUUNDARY LOURDINATES
        DATE
                                                                                                                                    NUMMALIZE THE CALLING COURDINATES
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                   ZOU HMI LL=HMIN+1RI-1.DO)*HCELL
AIML:L=ALMIN+1RJ-1.DO)*ALGELL
AFMCEL=AFMIN+1RK-1.DO)*APGELL
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                   MULT (XI, NCRDER, P1, TX)
MULT (XZ, NCRUER, P2, TY)
MULT (X3, NORDER, P3, TZ)
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                AAM-IX(NX)*IY(NY)*1Z(NZ)
(!!-6L'AAAA*CL!I)
(!|-6E+AAA*CE!II)
(.\-(\n^+ \aappa \appa 
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                        A)-(H-HMCELL)/HCELL
A-(ALAM-ALMCEL)/ALCELL
y-(APH)-APMCEL)/APCELL
    FINITE
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                          NY-17(11)
NY-17(11)
NY-17(11)
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IF THE TEST PUINT DUES NUT LIE WITHIN THE REGION MUDELED IN SECTION
    22/11/43
                                                         CALL PIMASS
LUSL=DLOS(ALAM)
SINL=DLOS(ALAM)
SINL=DSIN(ALAM)
COSP=DCUS(APHI)
COSP=DCUS(APHI)
SINL=SINL*CX + CUSP*SINL*GY + SINP*GZ
GU=COSP*COSL*GX + COSL*GY
GE=-SINP*CUSL*GX - SINF*SINL*GY + CCSP*GZ
Whit (0,1)
FORMAT(49H PUINT NET IN FINITE ELEMENT FIELD, FIMASS CALLED)
ISTITN=0
RETURN
  DATE = 79206
FINITE
RELEASE 2.0
                    ددون
FCRIKAN IV 61
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SM11H FROM COMMON XYZ

X,Y,Z--EAKIM-FIXED RECTANGULAR COCRDINATES OF PUINT AT WHICH GRAVITY
FROM CLAMEN MASPUS

FROM CLAMEN MASPUS

FROM CLAMEN MASPUS

FROM COMMENTED PRODUCTS OF THE GRAVITATIONAL CONSTANT AND THE

1000 PUINT MASSES (-1.6.), 6., 6., 1.6.)

PUSITS--PRECUMPUTED EARTH-FIXED X,Y,Z COORDINATES OF THE 1080 POINT CIVEN THE RELIANSULAR CCURDINATES, X, Y, Z, GF A POINT, THIS RUUTINE RETURNS THE CEMPCINENTS OF THE SKAVITY DISTURBANCE, DELGX, DELGY, AND DELGZ, USING THE POINT MASSES OF MASS MODEL SIG. TO COMMON XYZ LELGGEDFLOY, DELGZ--EAKTH-FIXED RECTANGULAR COMPONENTS OF THE GRAVITY DISTORBANCE FCR P. SOME POINT IN SPACE
DISTURBANCE IS TO BE EVALUATED
XI.YI.4.1.—EANIM-FIXED NECLANGULAR COUPDINATES OF THE ITH POINT MASS
PUSITS(1):1)=X1
FUSITS(1):2)=X1
FUSITS(1):2)=X1
FUSITS(1):2)=X1 -NHOO MODIFICATIONS BY JUHN L. JUNKINS, KEMI C. ENGELS, AND JOH DEPARIMENT DE ENGINEEKING SCIENCE AND MECHANICS SCHOOL OF ENGINEEKING VIRGINIA POLYTECHNIC INSTITUTE AND STATE UNIVERSITY ELACKSBURG, VIRGINIA 24061 JOHN L. JUNKINS AND JOHN SAUNDERS 79206 IMPLICIT REAL#8 (A-F.O-Z)
COMMUN /XYZ/X,Y,Z,DELGX,DELGY,DELGZ
COMMUN /MASPUS/PMVALS(1080),PUSITS(1080,3) OF LAST MULIFICATION -- APRIL 1,1979 SUBRCUTINE FIMASS 50 20 1=1,10EG 5EL67=0.00 0EL67=0.00 0EL64=0.00 MASSES JAJE PRCCESS CUPULS 1NPU1S 73 KELEASE FORTRAN IV G1 00000 2000 8000

FORTRAN IV G1

IF (PMVALS(I)) 10,20,10 CONTINUE DX=PUSITS(I,1)-X DY=PUSITS(I,2)-Y DZ=PLSITS(I,3)-Z UIST=CSUKT(UISTSU) TEMP=PMVALS(I)/DISTSU

DELGX=DELGX+(DX\*TEMP)
DELGY=DELGY+(UY\*TEMP)
DELGZ=UELGZ+(UZ\*TEMP)

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20 CUNTINUE

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ברי	:Icour,	15.10	26.62-									
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	10000		57-	1.6665	-0.1012	~~すつ・つー		5.5116		-5.3405	2.6969	1.1693
	103.35	_	57-	0.54.71	-4.3664	0.7126		2.4634	9	-5.5640	5.6361	-1.150
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SECURITY CLASSIFICATION OF THIS PAGE (When Date Entered)

REPORT DOCUMENTATION		READ INSTRUCTIONS BEFORE COMPLETING FORM
1. REPORT NUMBER	2. GOVT ACCESSION NO.	3. RECIPIENT'S CATALOG NUMBER
## TITLE (and Substitle) FINITE ELEMENT MODELS OF THE EAR'S FIELD PHASE IV	TH'S GRAVITY	5. TYPE OF REPORT & PERIOD COVERED  Contract Report 6. PERFORMING ORG. REPORT NUMBER
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7. Author(*) R. C. Engels, Assistant Professor J. L. Junkins, Professor, PI	r	DAAK-78-C-0072
<ol> <li>PERFORMING ORGANIZATION NAME AND ADDRESS Department of Engineering Science Virginia Polytechnic Institute Blacksburg, VA 24061</li> </ol>		10. PROGRAM ELEMENT, PROJECT, TASK AREA & WORK UNIT NUMBERS
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19. KEY WORDS (Continue on reverse elde if necessary as Gravity, Finite Element, Guidance	A BOT ARTS THE ARES	
Methods are developed and demo models for gravity anomalies. The efficient (by a factor of about 20) software listings and test cases.	onstrated for cons approach is show	structing piecewise local n to be computationally more